Part IV: PHYSICAL PROCESSES

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CHAPTER 1  Overview

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1.1 INTRODUCTION

Figure 1.1  Schematic diagram of the different physical processes represented in the IFS model.

The physical processes associated with radiative transfer, turbulent mixing, subgrid-scale orographic drag, moist convection, clouds and surface/soil processes have a strong impact on the large scale flow of the atmosphere. However, these mechanisms are often active at scales smaller than the horizontal grid size. Parametrization schemes are then necessary in order to properly describe the impact of these subgrid-scale mechanisms on the large scale flow of the atmosphere. In other words the ensemble effect of the subgrid-scale processes has to be formulated in terms of the resolved grid-scale variables. Furthermore, forecast weather parameters, such as two-metre temperature, precipitation and cloud cover, are computed by the physical parametrization part of the model.

This part (Part IV ‘Physical processes’) of the IFS documentation describes only the physical parametrization package. After all the explicit dynamical computations per time-step are performed, the physics parametrization package is called by the IFS. The physics computations are performed only in the vertical. The input information for the physics consists of the values of the mean prognostic variables (wind components, temperature, specific humidity, liquid/ice water content and cloud fraction), the provisional dynamical tendencies for the same variables
and various surface fields, both fixed and variable.

The time integration of the physics is based on the following:

1) it has to be compatible with the adiabatic part of the IFS;
2) the tendencies from the different physical processes are computed in separate routines;
3) as a general approach, the value of a prognostic variable is updated with the tendency from one process and the next process starts from this updated value, in what is usually referred to as the ‘method of fractional steps’ (details are different for different processes);
4) explicit schemes are used whenever possible, but if there are numerical stability problems the scheme is made as implicit as necessary.

The radiation scheme is described in Chapter 2 ‘Radiation’ and is the first process to be called in the physics. To save time in the rather expensive radiation computations, the full radiation part of the scheme is currently called every 3 hours. This is when the computation of the shortwave transmissivities and the longwave fluxes is performed, using the values of temperature, specific humidity, liquid/ice water content and cloud fraction at time-step \( t - \Delta t \), and a climatology for aerosols, CO\(_2\) and O\(_3\). The computation of the fluxes is not necessarily done at every grid-point but is only performed at sampled points, using a sampling algorithm that is latitude dependent. The results are then interpolated back to the original grid using a cubic interpolation algorithm. The shortwave fluxes are updated every time-step using synchronous values of the zenith angle. The radiation scheme takes into account cloud-radiation interactions in detail by using the values of cloud fraction and liquid/ice water content, at every level, from the prognostic cloud scheme. The radiation scheme produces tendencies of temperature.

The turbulent diffusion scheme is called just after radiation (Chapter 3 ‘Turbulent diffusion and interactions with the surface’). The surface fluxes are computed using Monin–Obukhov similarity theory. The computation of the upper-air turbulent fluxes is based on the \( K \)-diffusivity concept. Depending on the atmospheric stability different formulations for determining the \( K \)-coefficients are used: a \( K \)-profile closure for the unstable boundary layer and a \( Ri \)-number dependent closure for the stable boundary layer. Because of numerical stability problems the integration of the diffusion equation is performed in an implicit manner. In fact, it uses a so-called ‘more than implicit’ method, in which the ‘implicitness factor’ \( \alpha \) (which takes the value 0 in a fully explicit scheme and 1 in a fully implicit one) is set to 1.5. During the integration it uses the values of the prognostic variables at \( t - \Delta t \) to compute the \( K \)-coefficients but uses the tendencies updated by the dynamics and radiation on the right hand side of the discretized diffusion equation. The turbulent diffusion scheme also predicts the skin temperature and the apparent surface humidity. The turbulent diffusion scheme produces tendencies of temperature, specific humidity and wind components. It does not compute fluxes or tendencies of the cloud variables (liquid/ice water content and cloud fraction).

The moist convection scheme is described in Chapter 5 ‘Convection’. The scheme is based on the mass-flux approach and is divided in deep, mid-level and shallow convection. For deep convection the convective mass-flux is determined by assuming Convective Available Potential Energy (CAPE) is adjusted towards zero over a specified time-scale. For mid-level convection the cloud base mass-flux is directly related to the large scale vertical velocity. The intensity of shallow convection is estimated by assuming an equilibrium of moist static energy in the sub-cloud layer. The convection scheme provides tendencies of temperature, specific humidity and wind components.
In Chapter 6 ‘Clouds and large-scale precipitation’ the prognostic cloud scheme is described. It solves two prognostic equations for liquid/ice water content and cloud fraction. The cloud scheme represents the cloud formation by cumulus convection, the formation of boundary layer and stratiform clouds. The scheme also takes into account several important cloud processes like cloud-top entrainment, precipitation of water and ice and evaporation of precipitation. In the numerical integration of the equations the terms depending linearly on the values of liquid/ice water and cloud fraction are integrated analytically. The cloud scheme produces tendencies of all the prognostic variables.

The soil/surface scheme is described in Chapter 7 ‘Surface parametrization’. The scheme includes prognostic equations for temperature and moisture in four soil layers and snow mass. The soil equations use an implicit time integration scheme. An interception layer collects water from precipitation and dew fall. The evaporative fluxes consider separately the fractional contributions from snow cover, wet and dry vegetation and bare soil.

Chapter 8 ‘Methane oxidation’ describes a simple parametrization of the upper-stratospheric moisture source due to methane oxidation. A parametrization representing photolysis of vapour in the mesosphere is also included.

Chapter 9 ‘Ozone chemistry parametrization’ gives a brief description of the ozone parametrization and Chapter 10 ‘Climatological data’ describes the distributions of climatological fields.

1.2 OVERVIEW OF THE CODE

CALLPAR, the routine that controls the physical parametrization package, is called by CPGLAG that controls the grid-point calculations. CALLPAR calls the routines from the physics, the exception being the main radiation routine RADINT. RADINT controls the computation of the shortwave transmissivities and the longwave fluxes. RADINT is called via an interface routine RADDRV called by SCAN2MDM that is the multi-tasking interface to the computations in grid-point space (distributed memory version). RADINT is called outside CALLPAR because of the need to make the radiation space interpolation compatible with the distributed memory version of the IFS.

In CALLPAR the physics routines are called in the following order:

RARDERF: Computes radiative properties of the surface.
CLDP: Computes cloud parameters required for the post processing (e.g. total cloud cover).
RADHEATN: Computes the temperature tendencies and the downward radiation fluxes at the surface with updated (every time-step) values for the zenith angle.
VDFMAIN: Controls the computation of the vertical exchange of \(u, v, T\) and \(q\) by turbulence.
GWDRAG: Controls the computation of the tendencies for \(u, v, T\) due to the parametrization of subgrid-scale orographic drag.
CUCALLN: Interface to call CUMASTRN that controls the computation of the tendencies for \(u, v, T\) and \(q\) due to the parametrization of moist convective processes.
CLOUDSC: Controls the computation of tendencies for \(u, v, T, q, a\) and \(l\) due to the parametrization of the cloud processes.
SRFMAIN: Controls the soil/surface scheme.
METHOX: Computes tendencies for \(q\) due to methane oxidation and water vapour photolysis.
CHAPTER 2  Radiation

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2.1 RADIATIVE HEATING

The radiative heating rate is computed as the divergence of net radiation fluxes $\mathcal{F}$:
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\[
\left( \frac{\partial T}{\partial t} \right)_{rad} = \frac{g}{c_p} \frac{\partial f}{\partial p},
\]

(2.1)

where \( c_p \) is the specific heat at constant pressure of moist air

\[
c_p = c_{p_{dry}} \left\{ 1 + \left( c_{p_{vap}} - c_{p_{dry}} \right) q / c_{p_{dry}} \right\},
\]

and \( c_{p_{dry}} \) and \( c_{p_{vap}} \) are the specific heats at constant pressure of dry air and water vapour, respectively. Sections 2.2 and 2.3 describe the computation of the longwave and shortwave radiative fluxes respectively. The solution of the radiative transfer equation to obtain the fluxes is unfortunately very expensive, and we cannot afford to do it more than every 3 hours at every fourth grid point. The interpolation scheme used for obtaining the radiative fluxes at every grid point and every time step for the relevant instantaneous temperature profile and solar zenith angle is described in Section 2.4.

A description of the inputs, in particular the climatologically defined quantities of radiative importance is given in Section 2.5. Finally, an alphabetical list of the subroutines of the radiation scheme is given in Section 2.6.

### 2.2 LONGWAVE RADIATION

Since cycle 22r3, two longwave radiation schemes are available in the ECMWF model, the pre-cycle 22r3 by Morcrette (1991), and the current longwave radiation transfer scheme, the Rapid Radiation Transfer Model (RRTM).

The rate of atmospheric cooling by emission-absorption of longwave radiation is

\[
\frac{\partial T}{\partial t} = \frac{g}{c_p} \frac{\partial f_{LW}}{\partial p}
\]

(2.2)

where \( F_{LW} \) is the net longwave radiation flux (the subscript ‘LW’ is omitted in the remainder of this section).

Assuming a non-scattering atmosphere in local thermodynamic equilibrium, \( f \) is given by

\[
f = \int_{-1}^{1} \int_{0}^{\infty} dv \left\{ \mathcal{L}_v(p_{surf}, \mu) t_v(p_{surf}, p, \mu) + \int_{p'}^{\infty} \mathcal{L}_v(p', \mu) dt_v \right\}
\]

(2.3)

where \( \mathcal{L}_v(p, \mu) \) is the monochromatic radiance at wavenumber \( v \) at level \( p \), propagating in a direction \( \theta \) (the angle that this direction makes with the vertical), where \( \mu = \cos \theta \) and \( t_v(p, p' r) \) is the monochromatic transmission through a layer whose limits are at \( p \) and \( p' \) seen under the same angle \( \theta \), with \( r = \sec \theta \). The subscript ‘surf’ refers to the earth’s surface.

Subsections 2.2.1 to 2.2.4 describe the pre-cycle 22r3 scheme, and Subsection 2.2.5 describes the RRTM scheme in cycle 22r3.

#### 2.2.1 The pre-cycle 22r3 scheme

After separating the upward and downward components (indicated by superscripts + and –, respectively), and integrating by parts, we obtain the radiation transfer equation as it is actually estimated in the longwave part of the
radiation code

\[ f_v^+(p) = [B_v(T_{\text{surf}}) - B_v(T_{0\nu})]t_v(p_{\text{surf}}, p; r) + B_v(T(p)) + \int_{p_0}^{p} t_v(p, p'; r) dB_v \]

\[ f_v^-(p) = [B_v(T_{\infty}) - B_v(T_{\text{top}})]t_v(p, 0; r) + B_v(T(p)) + \int_{p_{\text{top}}}^{p} t_v(p', p; r) dB_v \]

(2.4)

where, taking benefit of the isotropic nature of the longwave radiation, the radiance \( L_v \) of (2.3) has been replaced by the Planck function \( B_v(T) \) in units of flux, W m\(^{-1}\) (here, and elsewhere, \( B_v \) is assumed to always includes the \( \pi \) factor). \( T_{\text{surf}} \) is the surface temperature, \( T_{0\nu} \) that of the air just above the surface, \( T(p) \) is the temperature at pressure-level \( p \), \( T_{\text{top}} \) that at the top of the atmospheric model. The transmission \( t_v \) is evaluated as the radiance transmission in a direction \( \theta \) to the vertical such that \( r = \sec \theta \) is the diffusivity factor (Elsasser, 1942). Such an approximation for the integration over the angle is usual in radiative transfer calculations, and tests on the validity of this approximation have been presented by Rodgers and Walshaw (1966) and Liu and Schmetz (1988) among others. The use of the diffusivity factor gives cooling rates within 2% of those obtained with a 4-point Gaussian quadrature.

### 2.2.2 Vertical integration

The integrals in (2.4) are evaluated numerically, after discretization over the vertical grid, considering the atmosphere as a pile of homogeneous layers. As the cooling rate is strongly dependent on local conditions of temperature and pressure, and energy is mainly exchanged with the layers adjacent to the level where fluxes are calculated, the contribution of the distant layers is simply computed using a trapezoidal rule integration, but the contribution of the adjacent layers is evaluated with a 2-point Gaussian quadrature, thus at the \( i \)th level,

\[ \int_{p_{i-2}}^{p_i} t_v(p, p'; r) dB_v = \]

\[ \sum_{l} dB_v(l)w_l t_v(p_l, p_l; r) + \frac{1}{2} \sum dB_v(j)[t_v(p_j, p_j; r) + t_v(p_j, p_{j-1}; r)] \]

(2.5)

where \( p_l \) is the pressure corresponding to the Gaussian root and \( w_l \) is the Gaussian weight. \( dB_v(j) \) and \( dB_v(l) \) are the Planck function gradients calculated between two interfaces, and between mid-layer and interface, respectively.

### 2.2.3 Spectral integration

The integration over wavenumber \( \nu \) is performed using a band emissivity method, as first discussed by Rodgers (1967). The longwave spectrum is divided into six spectral regions.

1) \( 0 \) – 350 cm\(^{-1}\) & \( 1450 \) – 1880 cm\(^{-1}\)
2) \( 500 \) – 800 cm\(^{-1}\)
3) \( 800 \) – 970 cm\(^{-1}\) & \( 1110 \) – 1250 cm\(^{-1}\)
4) \( 970 \) – 1110 cm\(^{-1}\)
corresponding to the centres of the rotation and vibration-rotation bands of H$_2$O, the 15 μm band of CO$_2$, the atmospheric window, the 9.6 μm band of O$_3$, the 25 μm “window” region, and the wings of the vibration-rotation band of H$_2$O, respectively. Over these spectral regions, band fluxes are evaluated with the help of band transmissivities precalculated from the narrow-band model of Morcrette and Fouquart (1985) - See Appendix of Morcrette et al. (1986) for details. Over these spectral regions, band fluxes are evaluated with the help of band transmissivities precalculated from the narrow-band model of Morcrette and Fouquart (1985) - See Appendix of Morcrette et al. (1986) for details.

Integration of (2.4) over wavenumber $\nu$ within the $k$th spectral region gives the upward and downward fluxes as

$$f_k^+(p) = \{B_k(T_{surf}) - B_k(T_0)\}t_{B_k}(r\nu(T_{surf}, p), T_{surf}(p, p')) + B_k(T_p)$$

$$+ \int_{p'}^p dB_k\{r\nu(p, p'), T_{surf}(p, p')\}dB_k$$

$$f_k^-(p) = \{B_k(T_0) - B_k(T_{surf})\}t_{B_k}(r\nu(p, 0), T_{surf}(p, 0)) - B_k(T_p)$$

$$- \int_{p'}^p dB_k\{r\nu(p', p), T_{surf}(p', p)\}dB_k$$

The formulation accounts for the different temperature dependencies involved in atmospheric flux calculations, namely that on $T_p$, the temperature at the level where fluxes are calculated, and that on $T_{surf}$, the temperature that governs the transmission through the temperature dependence of the intensity and half-widths of the lines absorbing in the concerned spectral region. The band transmissivities are non-isothermal accounting for the temperature dependence that arises from the wavenumber integration of the product of the monochromatic absorption and the Planck function. Two normalized band transmissivities are used for each absorber in a given spectral region: the first one for calculating the first right-hand-side term in (2.4), involving the boundaries; it corresponds to the weighted average of the transmission function by the Planck function

$$t_B(\nu, T_{surf}, T_0) = \int_{\nu_1}^{\nu_2} B_\nu(T_p)\nu d\nu$$

the second one for calculating the integral term in (2.4) is the weighted average of the transmission function by the derivative of the Planck function

$$t_B'(\nu, T_{surf}, T_0) = \int_{\nu_1}^{\nu_2} B_\nu(T_p)\nu d\nu$$
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\[
\int \{dB(T_p)/dT\} t_{dB} (\bar{\nu} p, T_u, T_u) \, d\nu
\]

\[
t_{dB} (\bar{\nu} p, T_p, T_u) = \frac{\int \{dB(T_p)/dT\} \, d\nu}{\int \{dB(T_p)/dT\} \, d\nu}
\]

(2.9)

where \( \bar{\nu} p \) is the pressure weighted amount of absorber.

The effect on absorption of the Doppler broadening of the lines (important only for pressure lower than 10 hPa) is included simply using the pressure correction method of Fels (1979). A finite line width (assumed to represent the Doppler half-width of the line) is retained under low pressure conditions where the pure Lorentz line width (proportional to pressure) would normally become negligible (Giorgetta and Morcrette, 1995).

In the scheme, the actual dependence on \( T_p \) is carried out explicitly in the Planck functions integrated over the spectral regions. Although normalized relative to \( B(T_p) \) or \( dB(T_p)/dT \), the transmissivities still depend on \( T_u \), both through Wien’s displacement of the maximum of the Planck function with temperature and through the temperature dependence of the absorption coefficients. For computational efficiency, the transmissivities have been developed into Pade approximants

\[
t(\bar{\nu} p, T_u) = \frac{\sum c_i \nu_i^{l/2}}{\sum d_j \nu_j^{l/2}}
\]

(2.10)

where \( \nu_{\text{eff}} = r(\bar{\nu} p) \Psi(T_u, \bar{\nu} p) \) is an effective amount of absorber which incorporates the diffusivity factor \( r \), the weighting of the absorber amount by pressure \( \bar{\nu} p \), and the temperature dependence of the absorption coefficients. The function \( \Psi(T_u, \bar{\nu} p) \) takes the form

\[
\Psi(T_u, \bar{\nu} p) = \exp[a(\bar{\nu} p)(T_u - 250) + b(\bar{\nu} p)(T_u - 250)^2]
\]

(2.11)

The temperature dependence due to Wien’s law is incorporated although there is no explicit variation of the coefficients \( c_i \) and \( d_j \) with temperature. These coefficients have been computed for temperatures between 187.5 and 312.5 K with a 12.5 K step, and transmissivities corresponding to the reference temperature the closest to the pressure weighted temperature \( T_u \) are actually used in the scheme.

2.2.4 The incorporation of the effects of clouds

The incorporation of the effects of clouds on the longwave fluxes follows the treatment discussed by Washington and Williamson (1977). Whatever the state of the cloudiness of the atmosphere, the scheme starts by calculating the fluxes corresponding to a clear-sky atmosphere and stores the terms of the energy exchange between the different levels (the integrals in (2.4)). Let \( f_0^+ (i) \) and \( f_0^- (i) \) be the upward and downward clear-sky fluxes. For any cloud layer actually present in the atmosphere, the scheme then evaluates the fluxes assuming a unique overcast cloud of emissivity unity. Let \( f_n^+ (i) \) and \( f_n^- (i) \) the upward and downward fluxes when such a cloud is present in the \( n \) th layer of the atmosphere. Downward fluxes above the cloud, and upward fluxes below the cloud, are assumed to be given by the clear-sky values.
\[ \mathcal{F}_n^+(i) = \mathcal{F}_0(i) \quad \text{for } i \leq n \]
\[ \mathcal{F}_n^-(i) = \mathcal{F}_0(i) \quad \text{for } i > n \]

Upward fluxes above the cloud (\( \mathcal{F}_n^+(k) \) for \( k \geq n + 1 \)) and downward fluxes below it (\( \mathcal{F}_n^-(k) \) for \( k > n \)) can be expressed with expressions similar to (2.5) provided the boundary terms are now replaced by terms corresponding to possible temperature discontinuities between the cloud and the surrounding air

\[ \mathcal{F}_n^+(k) = \{ \mathcal{F}_{\text{clld}}^+ - B(n + 1) \} t(p_k, p_{n+1}; r) + B(k) + \int_{p_n}^{p_k} t(p_k, p'; r) dB \]
\[ \mathcal{F}_n^-(k) = \{ \mathcal{F}_{\text{clld}}^- - B(n) \} t(p_k, p_n; r) + B(k) + \int_{p_n}^{p_k} t(p_k, p'; r) dB \]

where \( B(i) \) is now the total Planck function (integrated over the whole longwave spectrum) at level \( i \), and \( \mathcal{F}_{\text{clld}}^+ \) and \( \mathcal{F}_{\text{clld}}^- \) are the longwave fluxes at the upper and lower boundaries of the cloud. Terms under the integrals correspond to exchange of energy between layers in clear-sky atmosphere and have already been computed in the first step of the calculations. This step is repeated for all cloudy layers. The fluxes for the actual atmosphere (with semi-transparent, fractional and/or multi-layered clouds) are derived from a linear combination of the fluxes calculated in previous steps with some cloud overlap assumption in the case of clouds present in several layers. Let \( N' \) be the index of the layer containing the highest cloud, \( c_{\text{clld}}(i) \) the fractional cloud cover in layer \( i \), with \( c_{\text{clld}}(0) = 1 \) for the upward flux at the surface, and with \( c_{\text{clld}}(N + 1) = 1 \) and \( \mathcal{F}_{N+1}^- = \mathcal{F}_0^- \) to have the right boundary condition for downward fluxes above the highest cloud.

Whereas the maximum and random overlap assumptions are also available in the code (Morcrette and Fouquart, 1986), the maximum-random overlap assumption is operationally used in the ECMWF model, and the cloudy upward \( \mathcal{F}^+ \) and downward \( \mathcal{F}^- \) fluxes are obtained as

\[ \mathcal{F}(i) = \mathcal{F}_0(i) \quad \text{for } i = 1 \]
\[ \mathcal{F}(i) = C_{\text{clld}}(i - 1) \mathcal{F}_{i-1}^+(i) + \sum_{n=0}^{N-1} c_{\text{clld}}(n) \mathcal{F}_n^+(i) \prod_{l=n+1}^{i} (1 - c_{\text{clld}}(l)) \quad \text{for } 2 \leq i \leq N + 1 \]
\[ \mathcal{F}(i) = C_{\text{clld}}(N) \mathcal{F}_N^+(i) + \sum_{n=0}^{N-1} c_{\text{clld}}(n) \mathcal{F}_n^+(i) \prod_{l=n+1}^{N} (1 - c_{\text{clld}}(l)) \quad \text{for } i \geq N + 2 \]

In case of semi-transparent clouds, the fractional cloudiness entering the calculations is an effective cloud cover equal to the product of the emissivity due to the condensed water and the gases in the layer by the horizontal coverage of the cloud layer, with the emissivity, \( \varepsilon_{\text{clld}} \), related to the condensed water amount by

\[ \varepsilon_{\text{clld}} = 1 - \exp(-k_{\text{abs}} q_{\text{w},LWP}) \]

where \( k_{\text{abs}} \) is the condensed water mass absorption coefficient (in \( \text{m}^2\text{kg}^{-1} \)) following Smith and Shi (1992).
2.2.5 The Rapid Radiation Transfer Model (RRTM)

As stated in Mlawer et al. (1997), the objective in the development of RRTM has been to obtain an accuracy in the calculation of fluxes and heating rates consistent with the best line-by-line models. It utilizes the correlated-k method and shows its filiation to the Atmospheric and Environmental Research, Inc. (AER) line-by-line model (LBLRTM, Clough et al., 1989, 1992, Clough and Iacono, 1995) through its use of absorption coefficients for the relevant k-distributions derived from LBLRTM. Therefore the k-coefficients in RRTM include the effect of the CKD2.2 water vapour continuum (Clough et al., 1989).

The main point in the correlated-k method (Lacis and Oinas, 1991; Fu and Liou, 1992) is the mapping of the absorption coefficient \( k(\nu) \) from the spectral space (where it varies irregularly with wavenumber \( \nu \)) to the \( g \)-space (where \( g(k) \) is the probability distribution function, i.e. the fraction of the absorption coefficients in the set smaller than \( k \)). The effect of this reordering is a rearrangement of the sequence of terms in the integral over wavenumber in the radiative transfer equation (RTE), which makes it equivalent to what would be done for monochromatic radiation.

In the ECMWF model, no provision is presently taken for scattering in the longwave. Therefore, in order to get the downward radiance, the integration over the vertical dimension is simply done starting from the top of the atmosphere, going downward layer by layer. At the surface, the boundary condition (in terms of spectral emissivity, and potential reflection of downward radiance) is computed, then, in order to get the upward radiance, the integration over the vertical dimension is repeated, this from the surface upward.

The spectrally averaged radiance (between \( \nu_1 \) and \( \nu_2 \)) emerging from an atmospheric layer is

\[
\overline{R} = \frac{1}{(\nu_1 - \nu_2)} \int_{\nu_1}^{\nu_2} dv \left[ R_0(\nu) + \int_{t_r}^{1} [B(\nu, T(t'_\nu)) - R_0(\nu)] dt' \right]
\]

(2.16)

where \( R_0 \) is the incoming radiance to the layer, \( B(\nu, T) \) is the Planck function at wavenumber \( \nu \) and temperature \( T \), \( t_r \) is the transmittance for the layer optical path, and \( t'_\nu \) is the transmittance at a point along the optical path in the layer. Under the mapping \( \nu \rightarrow g \), this becomes

\[
\overline{R} = \int_{0}^{1} dg \left[ B_{\text{eff}}(g, T_g) + [R_0(g) - B_{\text{eff}}(g, T_g)] \exp\left[-k(g, P, T) \frac{\Delta z}{\cos \phi}\right] \right]
\]

(2.17)

where \( B_{\text{eff}}(g, T) \) is an effective Planck function for the layer that varies with the layer’s transmittance such as to ensure continuity of flux across layer boundaries for opaque conditions. The dependence of the transmittance is now written in terms of the absorption coefficient \( k(g, P, T) \) at layer pressure \( P \) and temperature \( T \), the absorber density \( \rho \), the vertical thickness of the layer \( \Delta z \), and the angle \( \phi \) of the optical path.

For a given spectral interval, the domain of the variable \( g \) is partitioned into subintervals (see Table 2.2, number of \( g \)-points), each corresponding to a limited range of \( k(g) \) values and for which a characteristic value \( \kappa_j \) of the absorption coefficient is chosen. These \( \kappa_j \) are then used to compute the outgoing radiance

\[
\overline{R} = \sum_j W_j \left[ B_{\text{eff}} + (R_0 - B_{\text{eff}}) \exp\left(-\kappa_j \frac{\Delta z}{\cos \phi}\right) \right]
\]

(2.18)

where \( W_j \) is the size of the sub-intervals (\( \sum W_j = 1 \)).

The accuracy of these absorption coefficients has been established by numerous and continuing high-resolution
validations of LBLRTM with spectroscopic measurements, in particular those from the Atmospheric Radiation Measurement program (ARM). Compared to the original RRTM (Mlawer et al., 1997), the version used at ECMWF has been slightly modified to account for cloud optical properties and surface emissivity defined for each of the 16 bands over which spectral fluxes are computed. For efficiency reason, the original number of \( g \)-points (256 = 16 \times 16) has been reduced to 140 (see Table 2.2). Other changes are the use of a diffusivity approximation (instead of the three-angle integration over the zenith angle used in the original scheme) to derive upward and downward fluxes from the radiances, and the modification of the original cloud random overlapping assumption to include (to the same degree of approximation as used in the operational SW scheme) a maximum-random overlapping of cloud layers. Given the monochromatic form of the RTE, the vertical integration is simply carried out one layer at a time from the top-of-the-atmosphere to the surface to get the downward fluxes. The downward fluxes at the surface are then used with the spectral surface emissivities and the surface temperature to get the upward long-wave fluxes in each of the 140 subintervals. Then the upward fluxes are obtained in a similar fashion from the surface to the ToA.

For the relevant spectral intervals of the RRTM schemes, ice cloud optical properties are derived from Ebert and Curry (1992), and water cloud optical properties from Fouquart (1987). Whereas in the operational scheme the cloud emissivity used to compute the effective cloud cover is defined over the whole LW spectrum from spectrally averaged mass absorption coefficients and the relevant cloud water and/or ice paths (following Smith and Shi, 1992), in RRTM, the cloud optical thickness is defined as a function of spectrally varying mass absorption coefficients and relevant cloud water and ice paths, and is used within the true cloudy fraction of the layer. Alternate sets of cloud optical properties are also available for RRTM, based on Savijarvi and Raisanen (1997) for liquid water clouds, and Fu et al. (1998) for ice clouds.

### 2.3 Shortwave Radiation

The rate of atmospheric heating by absorption and scattering of shortwave radiation is

\[
\frac{dT}{dt} = \frac{g}{c_p} \frac{\partial \mathcal{F}_{SW}}{\partial p} \quad (2.19)
\]

where \( \mathcal{F}_{SW} \) is the net total shortwave flux (the subscript SW will be omitted in the remainder of this section).

\[
\mathcal{F}(\delta) = \int_0^\infty \int_0^{2\pi} \int_{\mu=1}^{\mu=-1} L_v(\delta, \mu, \phi) d\mu \, d\phi 
\]

is the diffuse radiance at wavenumber \( v \), in a direction given by the azimuth angle, \( \phi \), and the zenith angle, \( \theta \), with \( \mu = \cos \theta \). In (2.20), we assume a plane parallel atmosphere, and the vertical coordinate is the optical depth \( \delta \), a convenient variable when the energy source is outside the medium

\[
\delta(p) = \int \beta_{v_{\text{ext}}}(p') \, dp'
\]

\( \beta_{v_{\text{ext}}}(p) \) is the extinction coefficient, equal to the sum of the scattering coefficient \( \beta_{v_{\text{sc}}} \) of the aerosol (or cloud particle absorption coefficient \( \beta_{v_{\text{abs}}} \)) and the purely molecular absorption coefficient \( k_v \). The diffuse radiance \( L_v \)
is governed by the radiation transfer equation

\[
\mu \frac{d \zeta_{\nu}(\delta, \mu, \phi)}{d \delta} = \zeta_{\nu}(\delta, \mu, \phi) - \frac{\sigma_{\nu}(\delta)}{4} P_{\nu}(\delta, \mu, \phi, \phi_{0}) \zeta_{\nu}^{0} \exp(-\delta/\mu_{s}) \]

(2.22)

\( I_{0} \) is the incident solar irradiance in the direction \( \mu_{0} = \cos \theta_{0} \), \( \sigma_{\nu} \) is the single scattering albedo (= \( \beta_{\nu}^{scc}/k_{\nu} \)) and \( \Phi(\delta, \mu, \phi, \mu', \phi') \) is the scattering phase function which defines the probability that radiation coming from direction \((\mu', \phi')\) is scattered in direction \((\mu, \phi)\). The shortwave part of the scheme, originally developed by Fouquart and Bonnel (1980) solves the radiation transfer equation and integrates the fluxes over the whole shortwave spectrum between 0.2 and 4 \( \mu \text{m} \). Upward and downward fluxes are obtained from the reflectance and transmittances of the layers, and the photon-path-distribution method allows to separate the parametrization of the scattering processes from that of the molecular absorption.

### 2.3.1 Spectral integration

Solar radiation is attenuated by absorbing gases, mainly water vapour, uniformly mixed gases (oxygen, carbon dioxide, methane, nitrous oxide) and ozone, and scattered by molecules (Rayleigh scattering), aerosols and cloud particles. Since scattering and molecular absorption occur simultaneously, the exact amount of absorber along the photon path length is unknown, and band models of the transmission function cannot be used directly as in longwave radiation transfer (see Section 2.2). The approach of the photon path distribution method is to calculate the probability \( \Pi(\nu) d\nu \) that a photon contributing to the flux \( F_{\text{cons}} \) in the conservative case (i.e., no absorption, \( \omega_{\nu} = 1, k_{\nu} = 0 \)) has encountered an absorber amount between \( \nu \) and \( \nu + d\nu \). With this distribution, the radiative flux at wavenumber \( \nu \) is related to \( F_{\text{cons}} \) by

\[
F_{\nu} = F_{\text{cons}} \int_{\nu}^{\infty} \Pi(\nu) \exp(-k_{\nu} \nu) d\nu \]  
(2.23)

and the flux averaged over the spectral interval \( \Delta \nu \) can then be calculated with the help of any band model of the transmission function \( t_{\Delta \nu} \)

\[
F = \frac{1}{\Delta \nu} \int_{\Delta \nu} F_{\nu} d\nu = F_{\text{cons}} \int_{\Delta \nu} \Pi(\nu) t_{\Delta \nu}(\nu) d\nu \]  
(2.24)

To find the distribution function \( \Pi(\nu) \), the scattering problem is solved first, by any method, for a set of arbitrarily fixed absorption coefficients \( k_{\nu} \), thus giving a set of simulated fluxes \( F_{\nu} \). An inverse Laplace transform is then performed on (2.23) (Fouquart, 1974). The main advantage of the method is that the actual distribution \( \Pi(\nu) \) is smooth enough that (2.23) gives accurate results even if \( \Pi(\nu) \) itself is not known accurately. In fact, \( \Pi(\nu) \) need not be calculated explicitly as the spectrally integrated fluxes are
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\[
\mathcal{F} = \mathcal{F}_{\text{cons}} \tau_{\Delta \nu} (\langle \mathcal{U} \rangle) \quad \text{in the limiting case of weak absorption}
\]

\[
\mathcal{F} = \mathcal{F}_{\text{cons}} \tau_{\Delta \nu} (\langle \mathcal{U}^2 \rangle) \quad \text{in the limiting case of strong absorption}
\]

where \( \langle \mathcal{U} \rangle = \int_0^\infty \Pi (\mathcal{U}) \mathcal{U} d\mathcal{U} \) and \( \langle \mathcal{U}^2 \rangle = \int_0^\infty \Pi (\mathcal{U}) \mathcal{U}^2 d\mathcal{U} \).

The atmospheric absorption in the water vapour bands is generally strong, and the scheme determines an effective absorber amount \( \mathcal{U}_e \) between \( \langle \mathcal{U} \rangle \) and \( \langle \mathcal{U}^2 \rangle \) derived from

\[
\mathcal{U}_e = \ln \left( \frac{\mathcal{F}_{\text{cons}}}{\mathcal{F}_k} \right) / k_e \quad (2.25)
\]

where \( k_e \) is an absorption coefficient chosen to approximate the spectrally averaged transmission of the clear sky atmosphere

\[
k_e = \frac{1}{\mathcal{U}_{\text{tot}}} \ln \left( \frac{\tau_{\Delta \nu} (\mathcal{U}_{\text{tot}} / \mu_0)}{\mu_0} \right) \quad (2.26)
\]

where \( \mathcal{U}_{\text{tot}} \) is the total amount of absorber in a vertical column and \( \mu_0 = \cos \theta_0 \). Once the effective absorber amounts of \( \text{H}_2\text{O} \) and uniformly mixed gases are found, the transmission functions are computed using Padé approximants

\[
t_{\Delta \nu} (\mathcal{U}) = \sum_{i=0}^{N} \frac{a_i \mathcal{U}^{i-1}}{\sum_{j=0}^{N} b_j \mathcal{U}^{j-1}} \quad (2.27)
\]

Absorption by ozone is also taken into account, but since ozone is located at low pressure levels for which molecular scattering is small and Mie scattering is negligible, interactions between scattering processes and ozone absorption are neglected. Transmission through ozone is computed using (2.24) where \( \mathcal{U}_{O_3} \) the amount of ozone is

\[
\mathcal{U}_{O_3}^d = M \int_0^3 d\mathcal{U}_{O_3} \quad \text{for the downward transmission of the direct solar beam}
\]

\[
\mathcal{U}_{O_3}^u = r \int_0^3 d\mathcal{U}_{O_3} + \mathcal{U}_{O_3}^d (p_{\text{surf}}) \quad \text{for the upward transmission of the diffuse radiation}
\]

\( r = 1.66 \) is the diffusivity factor (see Section 2.2), and \( M \) is the magnification factor (Rodgers, 1967) used instead of \( r \) to account for the sphericity of the atmosphere at very small solar elevations

\[
M = 35/\sqrt{\mu_0^2 + 1} \quad (2.28)
\]

To perform the spectral integration, it is convenient to discretize the solar spectrum into subintervals in which the surface reflectance, molecular absorption characteristics, and cloud optical properties can be considered as constants. One of the main causes for such a spectral variation is the sharp increase in the reflectivity of the vegetation in the near-infrared. Also, water vapour does not absorb below 0.69 \( \mu \text{m} \) nor do liquid water clouds. Till June 2000,
the ECMWF shortwave scheme considered only two spectral intervals, one for the visible (0.2 - 0.69 μm), one for the near-infrared (0.69-4.00 μm) parts of the solar spectrum. From June 2000 to April 2002, the near-infrared interval was sub-divided into three intervals (0.69 - 1.19 - 2.38 - 4.00 μm) to account better for the spectral variations of the cloud optical properties. Till April 2002, all the molecular absorption coefficients (for O₂, H₂O, uniformly mixed gases) were derived from statistical models of the transmission function using spectroscopic parameters derived from various versions of the HITRAN database (Rothman et al., 1986, 1992). In April 2002, following the recomputation of all the molecular absorption coefficients from an updated version of the shortwave line-by-line model of Dubuisson et al. (1996) using spectroscopic data from HAWKS (2000), the ultraviolet and visible part of the spectrum are now considered in three spectral intervals (0.20 - 0.25 - 0.69 μm) making the scheme having a total of six spectral intervals over which the aerosol and cloud optical properties are also defined. The cut-off at 0.69 μm allows the scheme to be more computational efficient, in as much as the interactions between gaseous absorption (by water vapour and uniformly mixed gases) and scattering processes are accounted for only in the near-infrared interval(s).

### 2.3.2 Vertical integration

Considering an atmosphere where a fraction \( C_{\text{cld}}^{\text{tot}} \) (as seen from the surface or the top of the atmosphere) is covered by clouds (the fraction \( C_{\text{cld}}^{\text{tot}} \) depends on which cloud-overlap assumption is assumed for the calculations), the final fluxes are given as a weighted average of the fluxes in the clear sky and in the cloudy fractions of the column

\[
F(j) = C_{\text{cld}}^{\text{tot}} F_{\text{cld}}(j) + \left(1 - C_{\text{cld}}^{\text{tot}} \right)F_{\text{clr}}
\]

where the subscripts ‘clr’ and ‘cld’ refer to the clear-sky and cloudy fractions of the layer, respectively. In contrast to the scheme of Geleyn and Hollingsworth (1979), the fluxes are not obtained through the solution of a system of linear equations in a matrix form. Rather, assuming an atmosphere divided into homogeneous layers, the upward and downward fluxes at a given layer interface \( j \) are given by

\[
F(j) = \prod_{k=j}^{N} R_{\text{bot}}(k) \hspace{1cm} (2.29)
\]

\[
F^-(j) = F^+(j) R_{\text{top}}(j-1)
\]

where \( R_{\text{top}}(j) \) and \( T_{\text{bot}}(j) \) are the reflectance at the top and the transmittance at the bottom of the \( j \)th layer. Computations of \( R_{\text{top}} \)'s start at the surface and work upward, whereas those of \( T_{\text{bot}} \)'s start at the top of the atmosphere and work downward. \( R_{\text{top}} \) and \( T_{\text{bot}} \) account for the presence of cloud in the layer

\[
R_{\text{top}} = C_{\text{cld}} R_{\text{cld}} + \left(1 - C_{\text{cld}} \right) R_{\text{clr}}
\]

\[
T_{\text{bot}} = C_{\text{cld}} T_{\text{cld}} + \left(1 - C_{\text{cld}} \right) T_{\text{clr}}
\]

where \( C_{\text{cld}} \) is the cloud fractional coverage of the layer within the cloudy fraction \( C_{\text{cld}}^{\text{tot}} \) of the column.

#### 2.3.2 (a) Cloudy fraction of the layer

\( R_{\text{cld}} \) and \( T_{\text{cld}} \) are the reflectance at the top and transmittance at the bottom of the cloudy fraction of the layer calculated with the Delta-Eddington approximation. Given \( \delta_{c} \), \( \delta_{a} \), and \( \delta_{g} \), the optical thicknesses for the cloud, the aerosol and the molecular absorption of the gases, respectively, and \( \left( = k_c U \right) \), and \( g_c \) and \( g_a \) the cloud and aerosol asymmetry factors, \( R_{\text{cld}} \) and \( T_{\text{cld}} \) are calculated as functions
of the total optical thickness of the layer

$$\delta = \delta_c + \delta_a + \delta_g$$

(2.31)

of the total single scattering albedo

$$\omega^* = \frac{\delta_c + \delta_a}{\delta_c + \delta_a + \delta_g}$$

(2.32)

of the total asymmetry factor

$$\varphi^* = \frac{\delta_c}{\delta_c + \delta_a} \varphi_c + \frac{\delta_a}{\delta_c + \delta_a} \varphi_a$$

(2.33)

of the reflectance $R_\omega$ of the underlying medium (surface or layers below the $j$th interface), and of the cosine of an effective solar zenith angle $\mu_{\text{eff}}(j)$ which accounts for the decrease of the direct solar beam and the corresponding increase of the diffuse part of the downward radiation by the upper scattering layers

$$\mu_{\text{eff}}(j) = \left[ 1 - \mathcal{C}_{\text{eff}}(j) \right] \mu + \mathcal{R}$$

(2.34)

with $\mathcal{C}_{\text{eff}}(j)$ the effective total cloudiness over level $j$

$$\mathcal{C}_{\text{eff}}(j) = 1 - \prod_{i=1}^{n} (1 - \mathcal{C}_{\text{cl}}(i) \mathcal{E}(i))$$

(2.35)

and

$$\mathcal{E}(i) = 1 - \exp \left[ -\frac{(1 - \mathcal{C}_{\text{cl}}(i) \mathcal{g}_{\varphi}(i))^2 \delta_c(i)}{\mu} \right]$$

(2.36)

$\delta_c(i)$, $\mathcal{C}_{\text{cl}}(i)$ and $\mathcal{g}_{\varphi}(i)$ are the optical thickness, single scattering albedo and asymmetry factor of the cloud in the $i$th layer, and $\mathcal{R}$ is the diffusivity factor. The scheme follows the Eddington approximation first proposed by Shettle and Weinman (1970), then modified by Joseph et al. (1976) to account more accurately for the large fraction of radiation directly transmitted in the forward scattering peak in case of highly asymmetric phase functions. Eddington’s approximation assumes that, in a scattering medium of optical thickness $\delta^*$, of single scattering albedo $\omega^*$, and of asymmetry factor $\varphi^*$, the radiance $\mathcal{L}$ entering (2.17) can be written as

$$\mathcal{L}(\delta, \mu) = \mathcal{L}_0(\delta) + \mu \mathcal{L}_1(\delta)$$

(2.37)

In that case, when the phase function is expanded as a series of associated Legendre functions, all terms of order greater than one vanish when (2.20) is integrated over $\mu$ and $\varphi$. The phase function is therefore given by

$$P(\Theta) = 1 + \beta_1(\Theta) \mu$$

where $\Theta$ is the angle between incident and scattered radiances. The integral in (2.20) thus becomes
where

\[
\theta = \frac{\beta_1}{3} = \frac{1}{2} \int_{-1}^{+1} P(\Theta) \mu \, d\mu
\]

is the asymmetry factor.

Using (2.38) in (2.20) after integrating over \( \mu \) and dividing by \( 2\pi \), we get

\[
\mu \frac{d}{d\delta} (\mathcal{L}_0 + \mu \mathcal{L}_1) = -(\mathcal{L}_0 + \mu \mathcal{L}_1) + \Theta(\mathcal{L}_0 + \theta \mathcal{L}_1) + 1/4 \Theta \mathcal{F}_0 \exp(-\delta/\mu_0)(1 + 3 \theta \mu_0 \mu)
\]

(2.39)

We obtain a pair of equations for \( \mathcal{L}_0 \) and \( \mathcal{L}_1 \) by integrating (2.39) over \( \mu \)

\[
\begin{align*}
\frac{d\mathcal{L}_0}{d\delta} &= -3(1 - \Theta) \mathcal{L}_0 + \frac{3}{4} \Theta \mathcal{F}_0 \exp(-\delta/\mu_0) \\
\frac{d\mathcal{L}_1}{d\delta} &= -(1 - \Theta \theta) \mathcal{L}_1 + \frac{3}{4} \Theta \theta \mu_0 \mathcal{F}_0 \exp(-\delta/\mu_0)
\end{align*}
\]

(2.40)

For the cloudy layer assumed non-conservative (\( \Theta < 1 \)), the solutions to (2.39) and (2.40), for \( 0 \leq \delta \leq \delta^* \), are

\[
\begin{align*}
\mathcal{L}_0(\delta) &= C_1 \exp(-K\delta) + C_2 \exp(+K\delta) - \alpha \exp(-\delta/\mu_0) \\
\mathcal{L}_1(\delta) &= P\{C_1 \exp(-K\delta) - C_2 \exp(+K\delta) - \beta \exp(-\delta/\mu_0)\}
\end{align*}
\]

(2.41)

where

\[
\begin{align*}
K &= \{3(1 - \Theta)(1 - \Theta \theta)\}^{1/2} \\
P &= \{3(1 - \Theta)/(1 - \Theta \theta)\}^{1/2} \\
\alpha &= 3 \Theta \mathcal{F}_0 \mu_0 \{1 + 3 \theta(1 - \Theta)\}/\{4(1 - K^2 \mu_0^2)\} \\
\beta &= 3 \Theta \mathcal{F}_0 \mu_0 \{1 + 3 \theta(1 - \Theta)\} \mu_0^2 /\{4(1 - K^2 \mu_0^2)\}
\end{align*}
\]

The two boundary conditions allow to solve the system for \( C_1 \) and \( C_2 \); the downward directed diffuse flux at the top of the atmosphere is zero, i.e.,

\[
\mathcal{F}(0) = \left[\mathcal{L}_0(0) + \frac{2}{3} \mathcal{L}_1(0)\right] = 0
\]

which translates into
The upward directed flux at the bottom of the layer is equal to the product of the downward directed diffuse and direct fluxes and the corresponding diffuse and direct reflectance (\( \mathcal{R}_d \) and \( \mathcal{R}_c \), respectively) of the underlying medium:

\[
\mathcal{F}^+(\delta^+)=\left\{ \mathcal{R}_d(\delta^+)-\frac{2}{3}\mathcal{L}_1(\delta^+) \right\}
\]

\[
=\mathcal{R}_c\left\{ \mathcal{L}_0(\delta^+)-\frac{2}{3}\mathcal{L}_1(\delta^+) \right\}+\mathcal{R}_d\mu_0\mathcal{F}_0\exp(-\delta^+\mu_0)
\]

which translates into

\[
\begin{align*}
\{1-\mathcal{R}_c-2(1+\mathcal{R}_c)P/3\}C_1\exp(-K\delta^+)+\{1-\mathcal{R}_c+2(1+\mathcal{R}_c)P/3\}C_2(\mathcal{R}_d\mu_0\mathcal{F}_0)exp(-\delta^+\mu_0)
\end{align*}
\]

In the Delta-Eddington approximation, the phase function is approximated by a Dirac delta function forward-scatter peak and a two-term expansion of the phase function:

\[
P(\theta) = 2f(1-\mu) + (1-f)(1 + 3g'\mu)
\]

where \( f \) is the fractional scattering into the forward peak and \( g' \) the asymmetry factor of the truncated phase function. As shown by Joseph et al. (1976), these parameters are

\[
f = \frac{g}{2}
g' = \frac{g}{g+1}
\]

The solution of the Eddington’s equations remains the same provided that the total optical thickness, single scattering albedo and asymmetry factor entering (2.39)–(2.43) take their transformed values:

\[
\begin{align*}
\delta^* &= (1+\sigma)f\delta^+ \\
\omega' &= \frac{(1-f)\sigma}{1-\sigma f}
\end{align*}
\]

Practically, the optical thickness, single scattering albedo, asymmetry factor and solar zenith angle entering (2.39)-(2.42) are \( \delta^* \), \( \sigma^* \), \( g' \) and \( \mu_{eff} \) defined in (2.33) and (2.34).

2.3.2 (b) Clear-sky fraction of the layers. In the clear-sky part of the atmosphere, the shortwave scheme accounts for scattering and absorption by molecules and aerosols. The following calculations are practically done twice, once for the clear-sky fraction (1 − \( C_{\text{cl}d} \)) of the atmospheric column with \( \mu \) equal to \( \mu_0 \), simply
modified for the effect of Rayleigh and aerosol scattering, the second time for the clear-sky fraction of each individual layer within the fraction $C_{clr}^{\mu}$ of the atmospheric column containing clouds, with $\mu$ equal to $\mu_{e}$.

As the optical thickness for both Rayleigh and aerosol scattering is small, $R_{clr}(j - 1)$ and $T_{clr}(j)$, the reflectance at the top and transmittance at the bottom of the $j$ th layer can be calculated using respectively a first and a second-order expansion of the analytical solutions of the two-stream equations similar to that of Coakley and Chylek (1975). For Rayleigh scattering, the optical thickness, single scattering albedo and asymmetry factor are respectively $\delta_{R}$, $\omega_{R} = 1$, and $g_{R} = 0$, so that

$$R_{R} = \frac{\delta_{R}}{2\mu + \delta_{R}}$$

$$T_{R} = \frac{2\mu}{(2\mu + \delta_{R})}$$

(2.46)

The optical thickness $\delta_{R}$ of an atmospheric layer is simply

$$\delta_{R} = \delta_{R}^{*} \left( \frac{p(j) - p(j - 1)}{p_{surf}} \right)$$

(2.47)

where $\delta_{R}^{*}$ is the Rayleigh optical thickness of the whole atmosphere parametrized as a function of the solar zenith angle (Deschamps et al., 1983)

For aerosol scattering and absorption, the optical thickness, single scattering albedo and asymmetry factor are respectively $\delta_{a}$, $\omega_{a}$, with $1 - \omega_{a} \approx 1$ and $g_{a}$, so that

$$\text{den} = 1 + \{1 - \omega_{a} + \text{back}(\mu_{e})\omega_{a}\}(\delta_{a} / \mu_{e})$$

$$+ (1 - \omega_{a})(1 - \omega_{a} + 2\text{back}(\mu_{e})\omega_{a})(\delta_{a} / \mu_{e}^{2})$$

$$R(\mu_{e}) = \frac{(\text{back}(\mu_{e})\omega_{a}\delta_{a}) / \mu_{e}}{\text{den}}$$

$$T(\mu_{e}) = 1 / \text{den}$$

(2.49)

where back($\mu_{e}$) = $(2 - 3\mu_{e}\delta_{a})$ is the backscattering factor.

Practically, $R_{clr}$ and $T_{clr}$ are computed using (2.49) and the combined effect of aerosol and Rayleigh scattering comes from using modified parameters corresponding to the addition of the two scatterers with provision for the highly asymmetric aerosol phase function through Delta-approximation of the forward scattering peak (as in (2.40)–(2.41))

$$\delta^{*} = \delta_{R} + \delta_{a}(1 - \omega_{a}g_{a}^{2})$$

$$g^{*} = \frac{g_{a}}{1 + g_{a}(\delta_{R} + \delta_{a})}$$

$$\omega^{*} = \frac{\delta_{R}}{\delta_{R} + \delta_{a} + \frac{\delta_{a}}{\delta_{R} + \delta_{a}}(1 - \omega_{a})}$$

(2.50)
As for their cloudy counterparts, R_{clr} and T_{clr} must account for the multiple reflections due to the layers underneath

\[ R_{clr} = R(\mu_e) + R(\mu_r)/(1 - R^\prime R) \]  \hspace{1cm} (2.51)

and R_\lambda is the reflectance of the underlying medium \( R_\lambda = R(\lambda - 1) \) and r is the diffusivity factor.

Since interactions between molecular absorption and Rayleigh and aerosol scattering are negligible, the radiative fluxes in a clear-sky atmosphere are simply those calculated from (2.27) and (2.45) attenuated by the gaseous transmissions (2.25).

### 2.3.3 Multiple reflections between layers

To deal properly with the multiple reflections between the surface and the cloud layers, it should be necessary to separate the contribution of each individual reflecting surface to the layer reflectance and transmittances in as much as each such surface gives rise to a particular distribution of absorber amount. In case of an atmosphere including N cloud layers, the reflected light above the highest cloud consists of photons directly reflected by the highest cloud without interaction with the underlying atmosphere, and of photons that have passed through this cloud layer and undergone at least one reflection on the underlying atmosphere. In fact, (2.22) should be written

\[ \mathcal{F} = \sum_{i=0}^{\infty} \mathcal{F}_c \int_0^\infty \tau_i(\tau) t_{\Delta \lambda}(\tau) d\nu \]  \hspace{1cm} (2.52)

where \( \mathcal{F}_c \) and \( \tau_i(\tau) \) are the conservative fluxes and the distributions of absorber amount corresponding to the different reflecting surfaces.

Fouquart and Bonnel (1980) have shown that a very good approximation to this problem is obtained by evaluating the reflectance and transmittance of each layer (using (2.39) and (2.45)) assuming successively a non-reflecting underlying medium (\( R_\lambda = 0 \)), then a reflecting underlying medium (\( R_\lambda \neq 0 \)). First calculations provide the contribution to reflectance and transmittance of those photons interacting only with the layer into consideration, whereas the second ones give the contribution of the photons with interactions also outside the layer itself.

From those two sets of layer reflectance and transmittances (\( \tau_{00}, \tau_{b0} \)) and (\( R_{b\neq}, T_{b\neq} \)) respectively, effective absorber amounts to be applied to computing the transmission functions for upward and downward fluxes are then derived using (2.23) and starting from the surface and working the formulas upward

\[ U_{c0} = \ln(\tau_{b0}/T_{bc})/k_e \]
\[ U_{c\neq} = \ln(\tau_{b\neq}/T_{bc})/k_e \]
\[ U_{e0} = \ln(R_{b0}/R_{bc})/k_e \]
\[ U_{e\neq} = \ln(R_{b\neq}/R_{bc})/k_e \]  \hspace{1cm} (2.53)

where \( R_{bc} \) and \( T_{bc} \) are the layer reflectance and transmittance corresponding to a conservative scattering medium.

Finally the upward and downward fluxes are obtained as

\[ \mathcal{F}(\lambda) = \mathcal{F}_0 \{ R_{e0} t_{\Delta \lambda}(U_{e0}) + (R_{b\neq} - R_{e0}) t_{\Delta \lambda}(U_{e\neq}) \} \]  \hspace{1cm} (2.54)
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2.3.4 Cloud shortwave optical properties

As seen in Sub-section 2.3.2 (a), the cloud radiative properties depend on three different parameters: the optical thickness $\delta_c$, the asymmetry factor $g_c$, and the single scattering albedo $\omega_c$.

Presently the cloud optical properties are derived from Fouquart (1987) for the water clouds, and Ebert and Curry (1992) for the ice clouds.

$\delta_c$ is related to the cloud liquid water amount $\varrho_{LWP}$ by

$$\delta_c = \frac{3\varrho_{LWP}}{2r_e}$$

where $r_e$ is the mean effective radius of the size distribution of the cloud water droplets. Presently $r_e$ is parametrized as a linear function of height from 10 µm at the surface to 45 µm at the top of the atmosphere, in an empirical attempt at dealing with the variation of water cloud type with height. Smaller water droplets are observed in low-level stratiform clouds whereas larger droplets are found in mid-level cumuliform water clouds.

In the two-, four-, and six- spectral interval versions of the shortwave radiation scheme, the optical properties of liquid water clouds are defined from Fouquart (1987) and those for ice clouds from Ebert and Curry (1992). Alternative optical properties are also available for liquid water clouds (Slingo, 1989) and ice clouds (Fu, 1996).

The effective radius of the liquid water cloud particles is computed from the cloud liquid water content using the diagnostic formulation of Martin et al. (1994) and specified concentrations of cloud concentration nuclei over land and ocean. For ice clouds, the effective dimension of the cloud particles is diagnosed from temperature using a revision of the formulation by Ou and Liou (1995).

2.4 Horizontal interpolation

As stated in the introduction, the cost of the radiation scheme described in the previous sections is prohibitive if it were used to compute the radiative fluxes at every time step and every grid point of the model.

In order to cut down the computing costs, the full radiation scheme is only used every 3 hours (“full radiation time steps”) at every fourth grid point, and a spatial and temporal interpolation thus provides the relevant interaction of the shortwave radiative fluxes with the solar zenith angle at every time step and every grid point.

To do so we define an effective transmissivity $\tau_c$ at each of the model level such that:

$$\mathcal{F}_s = \tau_c S_0$$

(2.56)

where $\mathcal{F}_s$ is the solar (shortwave) flux and $S_0$ is the solar flux at the top of the atmosphere. The values of $\mathcal{F}_s$ is kept constant between full radiation time steps and the net fluxes are recomputed at every time step, using (2.56) with the correct solar angle for every grid point.

The interpolation is done only in the zonal direction because of the strong meridional variation of the radiative processes and of the internal organisation of the model code. For each row, variables relevant to the input of the radiation calculation are transformed to a coarser subgrid via simple sampling. On the coarser grid so defined, fluxes are then evaluated as described in the previous sections. Finally, effective solar transmissivities and longwave
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2.5 INPUT TO THE RADIATION SCHEME

2.5.1 Model variables

Temperature values are needed at the boundaries of the layers, where the fluxes are computed. They are derived from the full level temperatures with a pressure weighted interpolation

\[
T_{h+1/2} = T_h \left( \frac{p_{k+1/2}(p_{k+1/2} - p_k)}{p_{k+1/2}(p_{k+1/2} - p_k)} \right) + T_h \left( \frac{p_{k+1/2}(p_{k+1/2} - p_k)}{p_{k+1/2}(p_{k+1/2} - p_k)} \right)
\]

(2.57)

At the bottom of the atmosphere, either the surface temperature or the temperature at 2 m is used, while at the top of the atmosphere the temperature is extrapolated from the first full level and second half level temperatures.

2.5.2 Clouds

Cloud fraction, and liquid/ice water content is provided in all layers by the cloud scheme.

2.5.3 Aerosols

Horizontal distributions for four climatological types of aerosols (oceanic, desert, urban, and stratospheric background) are defined from T5 spectral coefficients, with fixed vertical distributions following Tanre et al. (1984).

2.5.4 Carbon dioxide, ozone and trace gases

Carbon dioxide, methane, nitrous oxide, CFC-11 and CFC-12 have constant volume concentrations of 353 ppm, 1.72 ppm, 0.31 ppm, 280 ppt, and 484 ppt respectively (IPCC/SACC, 1990).

Two climatologies are available for the ozone distribution. In the first one (NOZOCL = 0), the ozone mixing ratio \( q_{O_3} \) depends on height, latitude, longitude and season. Its vertical distribution is assumed to be such that its integral from 0 to the pressure \( p \) is

\[
\int_0^p q_{O_3} dp = \frac{a}{1 + (b/p)^{3/2}}
\]

(2.58)

The constants \( a \) and \( b \) are related to the total amount of ozone and the height of its maximum mixing ratio. They are imposed in terms of a limited series of spherical harmonics (T10) for the geographical distribution and a Fourier series for the seasonal variation. The total amount of ozone was taken from London et al. (1976) and the altitude of the maximum concentration was derived from Wilcox and Belmont (1977). Plots of these values can be found in the Appendix. In the second climatology (NOZOCL = 1), the ozone mixing ratio \( q_{O_3} \) depends on height, latitude and month, and is taken from Fortuin and Langematz (1995).

2.5.5 Ground albedo and emissivity

The background land albedo, \( \alpha_{bg} \), is interpolated to the model grid from the monthly mean values of a snow-free albedo produced for the combined 1982–1990 years. The albedo for that dataset was computed using the method
of Sellers et al. (1996), but with new maps of soil reflectance, new values of vegetation reflectance and the biophysical parameters described in Los et al. (2000). More information on the original data and plots of the monthly mean albedo are shown in Chapter 10.

Spectral albedos for parallel and diffuse radiation are needed by the radiative code. In addition, the surface energy balance equation (see Chapter 3 on vertical diffusion) needs a spectrally integrated parallel+diffused albedo, specified for each independent surface functional unit, tile. The procedure is summarized in Table 2.1. Over open water, the surface albedo for direct parallel radiation is a fit to low-flying aircraft measurements over the ocean given by Taylor et al. (1996)

$$\alpha_{sp} = \frac{0.037}{1.1\mu_0^{1.4} + 0.15}$$

(2.59)

For sea ice, monthly values based on Ebert and Curry (1993) albedos for the Arctic Ocean are interpolated to the forecast time. The bare sea ice albedo value in Ebert and Curry is taken as a representative value for summer, and the dry snow albedo value is used for the winter months. Values for the Antarctic are shifted by six months. Separate values for visible and near-infrared spectral bands are used. The time-varying snow albedo ($\alpha_{sn}$, see Chapter 7), is used for the exposed snow tile only. Finally, the average of the diffuse and parallel albedos are spectrally integrated for each tile.

<table>
<thead>
<tr>
<th>Table 2.1</th>
<th>DIFFUSE AND PARALLEL ALBEDO AND WINDOW EMISSIVITY FOR EACH TILES</th>
</tr>
</thead>
<tbody>
<tr>
<td>Tile</td>
<td>1</td>
</tr>
<tr>
<td>Description</td>
<td>Open sea</td>
</tr>
<tr>
<td>Diffuse albedo</td>
<td>0.06</td>
</tr>
<tr>
<td>Parallel albedo</td>
<td>Taylor et al. (1996)</td>
</tr>
<tr>
<td>Window emissivity</td>
<td>0.99</td>
</tr>
</tbody>
</table>

The thermal emissivity of the surface outside the 800–1250 cm$^{-1}$ spectral region is assumed to be 0.99 everywhere. In the window region, the spectral emissivity is constant for open water, sea ice, the interception layer and exposed snow tiles. For low and high vegetation and for shaded snow the emissivity depends on the water content in the top soil layer. Emissivity decreases linearly from 0.96 for soils at or above field capacity to 0.93 for soils at or below permanent wilting point. The same formulation is used for bare ground, except for desert areas ($\alpha_{sb} > 0.3$), where a value of 0.93 is used independently of the soil water content. Finally, a broadband emissivity is obtained by convolution of the spectral emissivity and the Planck function at the skin temperature.

2.5.6 Solar zenith angle

Equations to compute the annual variation of the solar constant $I$, the solar declination $\delta$, and the difference between solar time and official time can be found in Paltridge and Platt (1976). These equations are used to give the cosine of the solar angle at the ground. Because of the curvature of the earth, the zenith angle is not quite constant along the path of a sun ray. Hence a correction is applied to $\mu_0$ to give an average $\mu_0$ for the atmosphere:
\[
\mu_0 = \frac{H}{a} \quad (2.60)
\]

where \(a\) is the earth radius and \(H\) is the atmospheric equivalent height. \(H/a\) is fixed at 0.001277.

2.6 The Radiation Code

Routine RADHEAT or RADHEATN (depending whether the diagnostic or prognostic cloud scheme is used) is called at every time step to compute the radiative fluxes and heating using the solar zenith angle computed in CPGLAG and emissivities and transmissivities (PEMTU, PTRSOL) computed at full radiation time steps in RADINT. The other routines are called either once at the beginning of the run (SUECRAD and below) or once per full radiation step at the first row (ECRADFR and below), or at every full radiation time step for all rows. In this section, we briefly describe the function of each routine.

<table>
<thead>
<tr>
<th>Spectral intervals cm(^{-1})</th>
<th>Number of g-points</th>
<th>Gases included</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td></td>
<td>Troposphere</td>
</tr>
<tr>
<td>10-250</td>
<td>8</td>
<td>H(_2)O</td>
</tr>
<tr>
<td>250-500</td>
<td>14</td>
<td>H(_2)O</td>
</tr>
<tr>
<td>500-630</td>
<td>16</td>
<td>H(_2)O, CO(_2)</td>
</tr>
<tr>
<td>630-700</td>
<td>14</td>
<td>H(_2)O, CO(_2)</td>
</tr>
<tr>
<td>700-820</td>
<td>16</td>
<td>H(_2)O, CO(_2), CCl(_4)</td>
</tr>
<tr>
<td>820-980</td>
<td>8</td>
<td>H(_2)O, CFC11, CFC12</td>
</tr>
<tr>
<td>980-1080</td>
<td>12</td>
<td>H(_2)O, O(_3)</td>
</tr>
<tr>
<td>1080-1180</td>
<td>8</td>
<td>H(_2)O, CFC12, CFC22</td>
</tr>
<tr>
<td>1180-1390</td>
<td>12</td>
<td>H(_2)O, CH(_4)</td>
</tr>
<tr>
<td>1390-1480</td>
<td>6</td>
<td>H(_2)O</td>
</tr>
<tr>
<td>1480-1800</td>
<td>8</td>
<td>H(_2)O</td>
</tr>
<tr>
<td>1800-2080</td>
<td>8</td>
<td>H(_2)O</td>
</tr>
<tr>
<td>2080-2250</td>
<td>4</td>
<td>H(_2)O, N(_2)O</td>
</tr>
<tr>
<td>2250-2380</td>
<td>2</td>
<td>CO(_2)</td>
</tr>
<tr>
<td>2380-2600</td>
<td>2</td>
<td>N(_2)O, CO(_2)</td>
</tr>
<tr>
<td>2600-3000</td>
<td>2</td>
<td>H(_2)O, CH(_4)</td>
</tr>
</tbody>
</table>

N.B.: CCl\(_4\) and CFC22 are presented not accounted for in the ECMWF model.

2.6.1 Set-up routines

- SUECRAD provides the interface with the user, via the namelist NAERAD. It defines the constants of Table 2.2 and sets the configuration for the radiative computations (from SUPHEC).
- ECRADFR modifies the frequency of full radiative computations (from CNT4).
Chapter 2 ‘Radiation’

- **SUAERL** and **SUAERSN** set up the longwave and shortwave radiative characteristics of the aerosols (from SUERCAD).
- **SUAERH** defines the geographical distribution of aerosols, in terms of spectral coefficients (from UPDTIER).
- **SUAERV** defines the globally averaged vertical distribution of the aerosols (from SUERCAD).
- **SUCLOP** sets up the longwave and shortwave radiative properties of the ice and water clouds (from SUERCAD).
- **SUECOZO** computes the Legendre coefficients for the ozone distribution according to the time of the year, using the Fourier coefficients defined in DATA statements (from UPDTIER).
- **SULWN** sets up the coefficients for the longwave radiative computations (from SUERCAD).
- **SURDI** sets up the concentrations of radiatively active gases and security parameters for the radiative computations (from SUERCAD).
- **SUSAT** sets up position and altitude of geostationary satellites in case of diagnostic simulation of radiances by the model radiation scheme (from SUERCAD).
- **SUSWN** sets up the coefficients for the shortwave radiative computations (from SUERCAD).
- **UPDTIER** updates the time for full radiative computations (from ECRADFR).
- The routines SUAERH, SUECOZO are called only once per full radiation step, at the first row.
- **SURRTAB** precomputes the array linking gaseous optical thickness and the transmission function (RRTM). (called from SUERCAD).
- **SURRTFTR** includes all coefficients related to the $g$-point configuration (RRTM). (called from SUERCAD).
- **SURRTPK** defines the limits of the spectral intervals, and the coefficients of the spectrally defined and spectrally integrated Planck functions (RRTM). (called from SUERCAD).
- **SURRTRF** defines the pressure and temperature reference profiles used for the tabulation of the absorption coefficients (RRTM). (called from SUERCAD).
- **RRTM_CMBGBn**, for each of the 16 spectral intervals, remaps the absorption coefficients from 16 to the final number of $g$-points (called from RRTM_INIT_140GP).
- **RRTM_INIT_140GP** performs the $g$-point reduction from 16 per band to a band-dependant number (column 2 in Table 1). It also computes the relative weighting for the new $g$-point combinations (called from SUERCAD).
- **RRTM_KGBn** contain the various absorption coefficients for all gases relevant to the different spectral bands.

2.6.2 Main routines

- **RADINT** is called by **RADDRV** to launch the full radiation computations. Zonal mean diagnostic of the temperature, clouds and albedo are computed. Temperature is vertically interpolated. Depending on the value of the variable NRINT an interpolation of all input variables to a coarser grid may be carried out. It may be necessary to subdivide the latitude belt in a few parts for the actual calculation of radiative fluxes because of storage space limitations. For this reason a loop over these parts follows. Inside this loop a call to routine **RADLSW** provides solar and thermal fluxes for a subset of points of that latitude row. These fluxes are converted into transmissivities and emissivities and after completion of the whole latitude circle they are transferred to the full grid when the calculations are carried out with the coarse resolution (NRINT>1).
- **RADLSW** is the driver routine of the solar and thermal fluxes by calling specialized routines **SW** for shortwave radiation and either **RRTM_RRTM_140GP** or **LW** for longwave radiation.
2.6.3 Specialized routines

- **RADSRF** is called from **RADPAR/CALLPAR** to compute surface albedo and emissivity. It computes the gridpoint diffuse and parallel spectral albedos and a spectrally integrated albedo (for postprocessing). It also computes the emissivity inside and outside the window region, and the spectrally integrated emissivity. Finally, it computes spectrally integrated tile albedos to be used by the surface energy balance routine (see Chapter 3 on vertical diffusion).
- **LW** organizes the longwave computation by calling in turn **LWU, LWBV, LWC**.
- **LWU** computes the effective absorber amounts including the pressure and temperature dependencies in the spectral intervals of the longwave radiation scheme.
- **LWBV** calls **LWB** and **LWV**
- **LWB** computes the Planck function with relation to temperature for all levels and spectral intervals.
- **LWV** organizes the vertical integration by calling **LWVN** which deals with the contribution to the flux of the layers adjacent to the level of computation of flux, **LWVD** which deals with the contribution from the more distant layers, and **LWVB** which computes the contribution of the boundary terms.
- **LWT** and **LWTM** compute the relevant transmission functions needed in **LWVN, LWVD**, and **LWVB**.
- **LWC** introduces the effect of clouds on the longwave fluxes.
- **SW** organizes the shortwave computation by calling in turn **SWU, SW1S**, and **SW2S**.
- **SWU** computes the effective absorber amounts including the pressure and temperature dependencies of the absorption.
- **SW1S** and **SW2S** deal with the shortwave radiation transfer in the two spectral intervals used to describe the solar spectrum. They both call **SWCLR**, which deals with the conservative scattering processes (Rayleigh) and the scattering / absorption by aerosols in the totally clear sky part of the atmospheric column, then **SWR** which deals with the same processes for the clear sky layers in an otherwise cloudy column, and **SWDE** which computes the reflectivity and transmissivity of a layer including non-conservative scatterers (cloud particles) with the Delta-Eddington approximation.
- **SWT** and **SWTT1**, computes the relevant transmission functions.
- **RRTM_RRTM_140GP** organizes the longwave computation by calling in turn, within a loop on the individual vertical columns, **RRTM_ECRT_140GP, RRTM_SETCOEF_140GP, RRTM_GASABS1A_140GP** and **RRTM_RTRN1A_140GP**.
- **RRTM_ECRT_140GP** defines the surface spectral emissivity, and the spectral aerosol thickness, and the layer absorber amounts and cloud quantities as used in **RRTM**.
- **RRTM_SETCOEF_140GP** computes the indices and fractions related to the pressure and temperature interpolations. It also calculates the values of the integrated Planck function for each spectral band at the level and layer temperatures.
- **RRTM_GASABS1A_140GP** launches the calculation of the spectrally defined optical thickness for gaseous absorption. It calls **RRTM_TAUMOLn**.
- **RRTM_RTRN1A_140GP** computes the downward then upward fluxes, using a diffusivity-type approximation for the angle integration. Cloud overlap is treated with a generalized maximum/random overlap method. Adjacent layers are treated with maximum overlap, non-adjacent cloud groups are treated with random overlap. For adjacent cloud layers, cloud information is carried from the previous two layers.
2.6.4 Heating rate computation

- **RADHEAT** or **RADHEATN**, depending whether the diagnostic or the prognostic cloud scheme is used, recomputes at each time step the net radiative fluxes from the layers' effective emissivity and transmissivity, using the actual temperature and solar zenith angle. It also computes the downward longwave and shortwave radiation at the surface.

**APPENDIX A  LIST OF SYMBOLS**

<table>
<thead>
<tr>
<th>Symbol</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>$B_\nu$</td>
<td>Planck function integrated over the half sphere with the factor involving $\pi$ absorbed: in units of flux ($W m^{-2}$)</td>
</tr>
<tr>
<td>$C_{cl}$</td>
<td>fractional cloud cover</td>
</tr>
<tr>
<td>$c_p$</td>
<td>specific heat at constant pressure of moist air</td>
</tr>
<tr>
<td>$c_{p,dry}$</td>
<td>specific heat at constant pressure of dry air</td>
</tr>
<tr>
<td>$c_{p,vap}$</td>
<td>specific heat at constant pressure of water vapour</td>
</tr>
<tr>
<td>$E_\nu$</td>
<td>incident solar radiance in the direction $\theta_0$</td>
</tr>
<tr>
<td>$f$</td>
<td>radiative flux</td>
</tr>
<tr>
<td>$f$</td>
<td>fractional scattering into the forward peak</td>
</tr>
<tr>
<td>$g$</td>
<td>acceleration of gravity</td>
</tr>
<tr>
<td>$g$</td>
<td>asymmetry factor for aerosol scattering</td>
</tr>
<tr>
<td>$k$</td>
<td>absorption coefficient</td>
</tr>
<tr>
<td>$L_\nu$</td>
<td>monochromatic radiance at wavenumber $\nu$</td>
</tr>
<tr>
<td>$M$</td>
<td>magnification factor ($= 35/\sqrt{(\mu_0^2 + 1)}$)</td>
</tr>
<tr>
<td>$m_{o_3}$</td>
<td>ozone mixing ratio</td>
</tr>
<tr>
<td>$P$</td>
<td>scattering phase function</td>
</tr>
<tr>
<td>$p$</td>
<td>pressure</td>
</tr>
<tr>
<td>$\Pi(\mu) d\mu$</td>
<td>probability of a photon encountering an absorber amount between $\mu$ and $\mu + d\mu$</td>
</tr>
<tr>
<td>$q$</td>
<td>specific humidity</td>
</tr>
<tr>
<td>$r$</td>
<td>diffusivity factor ($= \sec \theta$)</td>
</tr>
<tr>
<td>$r_\nu$</td>
<td>mean effective radius of cloud water droplets</td>
</tr>
<tr>
<td>$R_\nu$</td>
<td>reflectance</td>
</tr>
<tr>
<td>$S_0$</td>
<td>solar flux at the top of the atmosphere</td>
</tr>
<tr>
<td>$T$</td>
<td>transmittance</td>
</tr>
<tr>
<td>$T$</td>
<td>temperature</td>
</tr>
<tr>
<td>$t_\nu$</td>
<td>monochromatic transmission at wavenumber $\nu$</td>
</tr>
<tr>
<td>$\mu$</td>
<td>absorber amount</td>
</tr>
<tr>
<td>$\alpha$</td>
<td>surface albedo</td>
</tr>
<tr>
<td>$\beta_{abs}$</td>
<td>cloud particle absorption coefficient</td>
</tr>
<tr>
<td>$\beta_{ext}$</td>
<td>extinction coefficient</td>
</tr>
<tr>
<td>$\beta_{sca}$</td>
<td>scattering coefficient</td>
</tr>
<tr>
<td>$\delta_g$</td>
<td>molecular absorption of gases</td>
</tr>
<tr>
<td>$\delta$</td>
<td>optical depth</td>
</tr>
<tr>
<td>$\epsilon_{cl}$</td>
<td>cloud emissivity</td>
</tr>
<tr>
<td>$\mu$</td>
<td>= $\cos \theta$</td>
</tr>
<tr>
<td>$\nu$</td>
<td>wavenumber</td>
</tr>
</tbody>
</table>
\( \sigma_v \) single scattering albedo \( (= \beta^{\text{sc}}_v / k_v) \)
\( \Phi \) scattering phase function
\( \varphi \) azimuth angle
\( \theta \) zenith angle
\( \theta_0 \) direction of incident solar beam
\( \Theta \) angle between incident and scattered radiances
Part IV: PHYSICAL PROCESSES

CHAPTER 3  Turbulent diffusion and interactions with the surface

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   3.2.2 Stability functions
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   3.2.4 Roughness lengths
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   3.3.1 General
   3.3.2 The exchange coefficients
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3.5 The skin temperature
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3.8 Diagnostic computations for postprocessing
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3.1 INTRODUCTION

The parametrization scheme described in this chapter represents the turbulent transfer of heat, momentum and moisture between the surface and the lowest model level and the turbulent transport of the same quantities between model levels. The scheme computes the physical tendencies of the four prognostic variables ($u$, $v$, $T$ and $q$) due to the vertical exchange by turbulent (non-moist) processes. These tendencies are obtained as the difference between the results of an implicit time-step from $t$ to $t + 1$. All the diagnostic computations (such as the calculation of the exchange coefficients, etc.) are done at time $t$. The surface boundary condition is formulated separately for 8 different tiles: water, ice, wet skin, low vegetation, exposed snow, high vegetation, snow under vegetation, and bare soil. The different tiles have their own surface energy balance and their own skin temperature. In this version
of the IFS, the mixture of land and ocean tiles is still not used, i.e. a grid box is either 100% ocean (water + ice) or 100% land (tile 3 to 8). Details about tiles are given in Chapter 7.

The equation for the vertical diffusion of any conservative quantity $\psi$ is:

$$\frac{\partial \psi}{\partial t} = \frac{1}{\rho} \frac{\partial}{\partial z} \left( \rho K \frac{\partial \psi}{\partial z} \right) = \frac{1}{\rho} \frac{\partial J_\psi}{\partial z}$$  \hspace{1cm} (3.1)

The vertical turbulent flux $J_\psi$ (positive downwards) is written using a first-order turbulence closure, where $K_\psi$ is the exchange coefficient. The goal of the vertical diffusion parametrization is to define the exchange coefficients and then to solve equation (3.1) with the following boundary conditions:

$$K_\psi \frac{\partial \psi}{\partial z} = 0 \quad \text{at} \quad p = p_{\text{top}}$$

$$K_\psi \frac{\partial \psi}{\partial z} \to \sum_{i=1}^{N_T} F_i C_{i\psi} |U(z)| (\psi(z) - \psi_{\text{surf}}) \text{as} \ z \to 0$$  \hspace{1cm} (3.2)

where $p_{\text{top}}$ is the pressure at the top of the atmosphere. For heat and moisture the surface boundary condition is provided tile by tile and fluxes are averaged over the $N_T$ tiles, weighted by their fraction $F_i$. The transfer coefficient $C_{i\psi}$ at the lowest model level depends upon the static stability. The variable $\psi_{\text{surf}}$ represents the value of $\psi$ at the surface. For heat and moisture, 8 tiles are used (see Chapter 9). For wind, a single tile is used with a no slip condition at the surface.

The vertical diffusion process is applied to the two horizontal wind components, $u$ and $v$, the specific humidity $q$ and the dry static energy $s$, where

$$s = c_{p\text{dry}} (1 + Iq) T + gz = c_p T + \phi$$  \hspace{1cm} (3.3)

where $I = c_{p\text{vap}} / c_{p\text{dry}} - 1$ and $c_{p\text{dry}}$, $c_{p\text{vap}}$, and $c_p$ are the specific heats at constant pressure of dry air, water vapour and moist air, respectively, and $\phi$ is the geopotential.

The problem is simplified by assuming that $\phi$ remains constant with respect to time during the turbulent diffusion process (even if in reality $T$ variations would modify $z(p)$). Exchange coefficients (with the dimension of a pressure thickness) are then computed for momentum and for heat (sensible plus latent) (the subscripts ‘M’, ‘H’ and ‘Q’ are used to identify the exchange coefficient for momentum, heat and humidity), with different formulations for the stable and the unstable case (depending on the sign of a stability parameter, either the Obukhov length or the bulk Richardson number in the surface layer). The implicit linear equations for the fluxes of momentum, firstly for $u$ and $v$ and secondly for $s$ and $q$, are solved by a Gaussian-elimination/back-substitution method.

The surface boundary condition is applied between the downward scanning elimination and the upward scanning back substitution. It involves a no-slip condition for $u$ and $v$ and the tile-by-tile solution of the surface energy balance for the boundary condition of $s$ and $q$. The water tile is an exception as it ignores the surface energy balance and uses the specified SST and the saturation specific humidity as boundary conditions.

Finally, the tendency of the variable temperature is computed, modified by the effects of local dissipation (it is assumed that there is no storage of turbulence kinetic energy) and moisture diffusion on $c_p$. The tiled surface fluxes of heat and moisture are also computed for later use by the surface scheme.
3.2 THE SURFACE LAYER

The surface layer approximation is applied between the lowest model level (about 10 m above the surface in the 60-level model) and the surface and for each tile separately. It is assume that the turbulent fluxes are constant with height and equal to the surface values. They can be expressed, using Monin–Obukhov similarity theory, in terms of the gradients of wind, dry static energy and specific humidity, which are assumed to be proportional to universal gradient functions of a stability parameter:

\[
\begin{align*}
\frac{\kappa z \, \partial u}{u_+ \, \partial z} &= \Phi_M\left(\frac{z}{L}\right) \\
\frac{\kappa z \, \partial \theta}{s_+ \, \partial z} &= \Phi_H\left(\frac{z}{L}\right) \\
\frac{\kappa z \, \partial q}{q_+ \, \partial z} &= \Phi_Q\left(\frac{z}{L}\right)
\end{align*}
\]  \hspace{1cm} (3.4)

The scaling parameters \(u_+, s_+\) and \(q_+\) are expressed in terms of surface fluxes:

\[
\begin{align*}
\rho u_+^2 &= J_M \\
\rho u_+ s_+ &= J_s \\
\rho u_+ q_+ &= J_q
\end{align*}
\]  \hspace{1cm} (3.5)

The stability parameter \(L\) is the Obukhov length defined as

\[
L = -u_+^3 \left(\frac{\kappa g Q_{0v}}{T_I}\right)
\]

with

\[
Q_{0v} = \frac{u_+ s_+ - (e_{p,vap} - e_{p,dry}) T_I u_+ q_+}{e_p} + \varepsilon T_I u_+ q_+
\]  \hspace{1cm} (3.6)

\(Q_{0v}\) is the virtual temperature flux in the surface layer, \(\kappa\) is the Von Kármán constant (=0.4), \(T_I\) is a reference temperature taken as a near-surface temperature (the temperature of the lowest atmospheric level) and \(\varepsilon = (R_{vap} / R_{dry}) - 1\), where \(R_{vap}\) and \(R_{dry}\) are the gas constants for water vapour and dry air, respectively.

In the surface layer, the gradient functions (3.4) can be integrated to profiles:

\[
\begin{align*}
u &= \frac{\tau_x}{\kappa pu_+} \left\{ \log\left(\frac{z_l + z_{0M}}{z_{0M}}\right) - \Psi_M\left(\frac{z_l + z_{0M}}{L}\right) + \Psi_M\left(\frac{z_{0M}}{L}\right) \right\} \\
v &= \frac{\tau_y}{\kappa pu_+} \left\{ \log\left(\frac{z_l + z_{0M}}{z_{0H}}\right) - \Psi_H\left(\frac{z_l + z_{0M}}{L}\right) + \Psi_H\left(\frac{z_{0H}}{L}\right) \right\} \\
\frac{s - s_{surf}}{s_{surf}} &= \frac{J_s}{\kappa pu_+} \left\{ \log\left(\frac{z_l + z_{0M}}{z_{0H}}\right) - \Psi_H\left(\frac{z_l + z_{0M}}{L}\right) + \Psi_H\left(\frac{z_{0H}}{L}\right) \right\} \\
\frac{q - q_{surf}}{q_{surf}} &= \frac{J_q}{\kappa pu_+} \left\{ \log\left(\frac{z_l + z_{0M}}{z_{0Q}}\right) - \Psi_Q\left(\frac{z_l + z_{0M}}{L}\right) + \Psi_Q\left(\frac{z_{0Q}}{L}\right) \right\}
\end{align*}
\]  \hspace{1cm} (3.7-3.10)
\[ z_{0M}, z_{0H}, \text{ and } z_{0Q} \text{ are the roughness lengths for momentum, heat and moisture. The stability profile functions } \Psi \text{ are derived from the gradient functions (3.4) with the help of the relationship } \Phi = 1 - \zeta (\partial \Psi / \partial \zeta). \text{ These profiles are used for the surface atmosphere interaction as explained in the following sections and also for the interpolation between the lowest model level and the surface (postprocessing of 10 m wind and 2 m temperature and moisture).} \]

In extremely stable situations, i.e. for very small positive \( L \), the ratio \( z / L \) is large, resulting in unrealistic profile shapes with standard stability functions. Therefore the ratio \( z / L \) is limited to 5 by defining a height \( h \) such that \( h / L = 5 \). If \( z < h \), then the profile functions described above, are used up to \( z = h \) and the profiles are assumed to be uniform above that. This modification of the profiles for exceptionally stable situations (no wind) is applied to the surface transfer formulation as well as to the interpolation for postprocessing.

### 3.2.1 Surface fluxes

Surface fluxes for heat and moisture are computed separately for the different tiles, so most of the surface layer computations loop over the tile index. Here a general description is given of the aerodynamic aspects of the transfer between the surface and the lowest model level. The description of the individual tiles can be found in Chapter 7.

Assuming that the first model level above the surface is located in the surface boundary layer at a specified height \( z_l \), the gradient functions (3.4) can be integrated to profiles for wind, dry static energy and specific humidity. The surface fluxes are expressed in terms of differences between parameters at level \( z_l \) and surface quantities (identified by the subscript ‘surf’; the tile index has been omitted in this general description).

\[ J_M = \rho C_M |U_l|^2 \\
J_s = \rho C_H |U_l|(s_l - s_{surf}) \]

\[ J_q = \rho C_Q |U_l|(\alpha_q - \alpha_{surf} q_{surf}) \]

where \( q_{surf} = q_{sa}(T_{surf}) \), \( \alpha_q \) and \( \alpha_{surf} \) are provided by the land scheme, \( s_{surf} = c_{po}(1 + \delta q_{sa})T_{surf} \), and \( q_{sa} \) is the apparent surface humidity also provided by the land surface scheme (the humidity equation simplifies over water where \( \alpha_q = 1 \), \( \alpha_{surf} = 1 \) and \( q_{sa} = q_{surf} \)).

The transfer coefficients can be expressed as follows

\[ C_M = \frac{\kappa^2}{\log (z_l / z_{0M}) - \Psi_M (z_l / z_{0M}) + \Psi_M (z_{0M} / L)^2} \]

\[ C_H = \frac{\kappa^2}{\log (z_l / z_{0M}) - \Psi_M (z_l / z_{0M}) + \Psi_M (z_{0M} / L)^2 \log (z_l / z_{0H}) - \Psi_H (z_l / z_{0H}) + \Psi_H (z_{0H} / L)^2} \]

\[ C_Q = \frac{\kappa^2}{\log (z_l / z_{0M}) - \Psi_M (z_l / z_{0M}) + \Psi_M (z_{0M} / L)^2 \log (z_l / z_{0Q}) - \Psi_Q (z_l / z_{0Q}) + \Psi_Q (z_{0Q} / L)^2} \]

The wind speed \(|U_l|\) is expressed as
with $w_*$ the free convection velocity scale defined by

$$w_* = \left( z_i \frac{g}{T_l} Q_{ov} \right)^{1/3}$$

The parameter $z_i$ is a scale height of the boundary layer depth and is set to constant value of 1000 m, since only the order of magnitude matters. The additional term in equation (3.15) represents the near surface wind induced by large eddies in the free-convection regime. When the surface is heated, this term guarantees a finite surface wind-forcing in the transfer law even for vanishing $u_l$ and $v_l$, and prevents $[U_l]^2$ and $L$ from becoming zero. Beljaars (1994) showed that this empirical term, when added into the standard Monin–Obukhov scaling, is in agreement with scaling laws for free convection. When used with the roughness lengths defined below, it provides a good fit to observational data, both over land and over sea.

### 3.2.2 Stability functions

The empirical forms of the dimensionless gradient functions $\Phi$ (equations (3.4)) have been deduced from field experiments over homogeneous terrain.

(a) In unstable conditions, $(\zeta = z/L < 0)$, the gradient functions proposed by Dyer and Hicks are used (Dyer, 1974; Hogström, 1988):

$$\Phi_M(\zeta) = (1 - 16\zeta)^{-1/4}$$

$$\Phi_H(\zeta) = \Phi_Q(\zeta) = (1 - 16\zeta)^{-1/2}$$

(3.17)

These functions can be integrated to the universal profile stability functions, $\Psi$, (Paulson, 1970):

$$\Psi_M(\zeta) = \frac{\pi}{2} - 2 \tan^{-1}(x) + \log \left( \frac{1 + x^2}{8} \right) \frac{1 + x^2}{8}$$

$$\Psi_H(\zeta) = \Psi_Q(\zeta) = 2 \log \left( \frac{1 + x^2}{2} \right)$$

with $x = (1 - 16\zeta)^{1/4}$. The $\Psi$-functions are used in the surface layer and the $\Phi$-functions for unstable stratification are used above the surface layer for local closure.

(b) For stable conditions, $(\zeta = z/L > 0)$, the code contains gradient function $\Phi_M$ as documented by Hogström (1988), and $\Phi_H$ as derived from the Ellison and Turner relation for the ratio $\Phi_M/\Phi_H$:

$$\Phi_M(\zeta) = 1 + 5\zeta$$

$$\Phi_H(\zeta) = \Phi_Q(\zeta) = (1 + 4\zeta)^2$$

(3.19)

These functions were meant to be used for local closure above the surface layer, but are not used at all in the current model version, because Richardson number dependent functions are used instead (see section on exchange coefficients above the surface layer).
The stable profile functions as used in the surface layer, are assumed to have the empirical forms proposed by Holtslag and De Bruin (1988), with a modification to allow for the effects of a critical flux Richardson number for large $\zeta$:

$$\Psi_M(\zeta) = -b\left(\zeta - \frac{c}{d}\right)\exp(-d\zeta) - a\zeta - \frac{bc}{d}$$

$$\Psi_H(\zeta) = \Psi_Q(\zeta) = -b\left(\zeta - \frac{c}{d}\right)\exp(-d\zeta) - \left(1 + \frac{2}{3}a\zeta\right)^{1.5} - \frac{bc}{d} + 1$$

(3.20)

where $a = 1$, $b = 2/3$, $c = 5$, and $d = 0.35$.

### 3.2.3 Computation of the Obukhov length

The transfer coefficients needed for the surface fluxes require the estimation of stability parameter $\zeta$, itself a function of the surface fluxes. Therefore, an implicit equation, relating $\zeta$ to bulk Richardson number $Ri_{\text{bulk}}$, is solved:

$$Ri_{\text{bulk}} = \frac{\zeta}{\left[\log\left(\frac{z_l+z_{0M}}{z_{0H}}\right) - \Psi_H\left(\frac{z_l+z_{0M}}{L}\right) + \Psi_H\left(\frac{z_{0H}}{L}\right)\right]}
\left[\log\left(\frac{z_l+z_{0M}}{z_{0M}}\right) - \Psi_M\left(\frac{z_l+z_{0M}}{L}\right) + \Psi_M\left(\frac{z_{0M}}{L}\right)\right]^{-1}$$

(3.21)

with

$$Ri_{\text{bulk}} = \left(\frac{g}{\theta_l}\right)\frac{z_l(\theta_{vl} - \theta_{v\text{surf}})}{|U_l|^2}$$

(3.22)

where $\theta_{vl}$ and $\theta_{v\text{surf}}$ are the virtual potential temperatures at level $z_l$ and at the surface, and $\theta_{vl}$ is a virtual potential temperature within the surface layer. Equation (3.22) can be expressed in terms of dry static energy:

$$Ri_{\text{bulk}} = \frac{g^2 z_l}{|U_l|^2}\left[\frac{2(s_l - s_{\text{surf}})}{(s_l + s_{\text{surf}} - \phi_l)} - (d - e)(q_l - q_{\text{surf}})\right]$$

(3.23)

Knowing $Ri_{\text{bulk}}$ at time $t$, a first guess of the Obukhov length is made from fluxes computed at the previous time step. Equation (3.21) is solved numerically using the Newton iteration method to retrieve $\zeta$.

In contrast to the previous formulation used in the model (Louis et al., 1982), the present scheme allows a consistent treatment of different roughness lengths for momentum, heat and moisture. The revised stability functions also reduce diffusion in stable situations resulting in more shallow stable boundary layers.

### 3.2.4 Roughness lengths

The integration constants $z_{0M}$, $z_{0H}$ and $z_{0Q}$, in the equations for the transfer coefficients $C_M$, $C_H$ and $C_Q$, (equations (3.12)–(3.14)) are called roughness lengths because they are related to the small scale inhomogeneities of the surface that determine the air–surface transfer.

- Over land, roughness lengths are assumed to be fixed climatological fields as described in Chapter 9. They are derived from land-use maps, with an extra contribution dependent on the variance of subgrid orography.
• Over sea, the specification of surface roughness lengths is particularly important. Because of the fixed boundary conditions for temperature and moisture the sea is, in principle, an infinite source of energy to the model. The surface roughness lengths are expressed by (Beljaars, 1994):

\[
\begin{align*}
 z_{0M} &= \alpha_M \frac{v}{u_*} + \alpha_{Ch} \frac{u_*^2}{g} \\
 z_{0H} &= \alpha_H \frac{v}{u_*} \\
 z_{0Q} &= \alpha_Q \frac{v}{u_*}
\end{align*}
\] (3.24)

These expressions account for both low and high wind regimes:

- At low wind speed the sea surface becomes aerodynamically smooth and the sea surface roughness length scales with the kinematic viscosity \( v(= 1.5 \cdot 10^{-5} \text{m}^2\text{s}^{-1}) \).

- At high wind speed the Charnock relation is used. The chosen constants are \( \alpha_M = 0.11 \), \( \alpha_H = 0.40 \), and \( \alpha_Q = 0.62 \) (Brutsaert, 1982). The Charnock coefficient, \( \alpha_{Ch} \), is set equal to 0.018 for the uncoupled model, and is provided by the wave model in coupled mode.

The smooth-surface parametrization is retained in high wind speed regimes for heat and moisture because observations indicate that the transfer coefficients for heat and moisture have very little wind-speed dependence above 4 m s\(^{-1}\) (Miller et al., 1992; Godfrey and Beljaars, 1991). In Eqs. (3.24), friction velocity \( u_* \), is calculated from

\[
u_* = C_M^{1/2} \left( u_l^2 + v_l^2 + w_*^2 \right)^{1/2}
\] (3.25)

with \( w_* \) from equation (3.16) using fluxes from the previous time step.

### 3.3 THE EXCHANGE COEFFICIENTS ABOVE THE SURFACE LAYER

#### 3.3.1 General

A first order closure specifies the turbulent flux of a given quantity \( \psi \) at a given model level proportional to the vertical gradient of that quantity:

\[
J_\psi = \rho K_\psi \frac{\partial \psi}{\partial z}
\] (3.26)

The exchange coefficients \( K_\psi \) are estimated at half model levels. The computation of the exchange coefficients depends on the stability regimes (locally and at the surface) and on the vertical location above the surface. Fig. 3.1 summarizes the various areas where each scheme (non-local diffusion from Troen and Mahrt, local diffusion dependent on the Richardson number, local diffusion with Monin–Obukhov functions) is applied. First, the local Richardson number is computed in each vertical layer:
Given the value of $\text{Ri}$, in stable local conditions the stability parameter $\zeta = z/L$ is deduced from precomputed tables giving $\zeta = \zeta(\text{Ri})$. A cubic spline interpolation is performed (Press et al. 1992, pp107-111). In unstable local conditions, we simply set $\zeta = \text{Ri}$.

### 3.3.2 The exchange coefficients

#### 3.3.2 (a) Turbulence length scale.

The mixing lengths $l_M = l_H = \kappa z$ used in the surface layer are bounded in the outer layer by introducing asymptotic length scales $\lambda_M \beta$ and $\lambda_H \beta$ (Blackadar, 1962)

$$
\frac{1}{l_M} = \frac{1}{\kappa z} + \frac{1}{\lambda_M \beta} \quad (3.28)
$$

$$
\frac{1}{l_H} = \frac{1}{\kappa z} + \frac{1}{\lambda_H \beta}
$$

The underlying idea is that vertical extent of the boundary layer limits the turbulence length scale. Since the results in the boundary layer are not very sensitive to the exact value of the asymptotic length scales, these parameters are chosen to be constants:

$$
\lambda_H = \lambda_M = 150 \text{ m}.
$$

Parameter $\beta$ is 1 in the boundary layer but reduces the length scales above the boundary layer in order to prevent excessive mixing to occur in and around the jet stream. The following expression is used

$$
\beta = \beta_o + \frac{(1 - \beta_o)}{1 + \left(\frac{z}{H_{\text{min}}}\right)^2}
$$

where $\beta_o = 0.2$ and $H_{\text{min}} = 4000 \text{ m}$.

#### 3.3.2 (b) $M$–$O$ similarity with $\text{Ri} < 0$ (Area 1 in Fig. 3.1).

In this regime, the exchange coefficients $K_\psi$ are based on local similarity (Nieuwstadt, 1984) stating that the expressions of the surface layer similarity can be used in the outer layer (strictly speaking only valid for stable conditions):
Here it is used for the unstable regime above the boundary layer, basically to provide strong vertical mixing in statically unstable situations.

3.3.2 (c) Revised Louis scheme for Ri > 0 (Area 1 in Fig. 3.1).

The use of Eq. (3.31) to define the exchange coefficients in the stable regime was found to be detrimental to the scores of the model (Beljaars, 1995) because of insufficient turbulent exchange in the lower troposphere. Therefore a revised version of the Louis scheme is used (Beljaars and Viterbo, 1999; Viterbo et al., 1999):

\[
K_M = \frac{i_M^2}{\Phi_M^2} \frac{\partial U}{\partial z}
\]

\[
K_H = \frac{i_H^2}{\Phi_M \Phi_H} \frac{\partial U}{\partial z}
\]

The functional dependencies of \( f_M \) and \( f_H \) with \( \text{Ri} \) are:

\[
f_M(\text{Ri}) = \frac{1}{1 + 2b \text{Ri}(1 + d \text{Ri})^{-1/2}}
\]

\[
f_H(\text{Ri}) = \frac{1}{1 + 2b \text{Ri}(1 + d \text{Ri})^{1/2}}
\]

with \( b = 5 \) and \( d = 1 \) (these functions are revised versions of the Louis et al., 1982 functions and were introduced in September 1995 in order to enhance turbulent transport in stable layers, see Viterbo et al., 1999).

3.3.2 (d) Unstable at the surface (Area 2 in Fig. 3.1). In unstable surface conditions (\( Q_{0v} < 0 \)), the exchange coefficients are expressed as integral profiles for the entire convective mixed layer. This K-profile closure is based on the form proposed by Troen and Mahrt (1986). This approach is more suitable than the local diffusion one when the length scale of the largest transporting turbulent eddies have a similar size as the boundary layer height itself (unstable and convective conditions). It also allows for an explicit entrainment parametrization in the capping inversion (Beljaars and Viterbo, 1999).

First a characteristic turbulent velocity scale \( w_{\text{turb}} \) is computed:

\[
w_{\text{turb}} = (u^*_t + 0.6w^*_{t})^{1/3}
\]

The velocities \( u^*_t \) and \( w^*_{t} \) are defined by equations (3.25) and (3.16) respectively.

Since the most energetic transporting scales of turbulent motion in the convective boundary layer are thermals, their strength is defined by a temperature excess with respect to their surrounding. The virtual dry static energy excess is written as
The profile is scanned to find the first virtual dry static energy minimum above the surface \( s_{vmin} \) at level \( z_{kmin} \).
The profile is further scanned to find mixed-layer depth, \( z_i \), defined in terms of the first level \( k \) where \( s_k > s_{vmin} + s_{turb} \), and \( z_i = 0.5(z_k + z_{k+1}) \).

- **Area 2.1 in Fig. 3.1**. In the surface layer above the first atmospheric level, \( z_l < z < 0.1z_i \), the exchange coefficients are prescribed as follows:

\[
K_M = kzw_{turb}\left(1 - \frac{z}{z_i}\right)^2 \frac{1}{\Phi_M} \\
K_H = kzw_{turb}\left(1 - \frac{z}{z_i}\right)^2 \frac{1}{\Phi_H}
\]  

(3.36)

- **Area 2.2 in Fig. 3.1**. In the unstable outer layer ( \( 0.1z_i < z \leq z_i \) ), similar expressions are used:

\[
K_M = kzw_{turb}\left(1 - \frac{z}{z_i}\right)^2 \\
K_H = kzw_{turb}\left(1 - \frac{z}{z_i}\right)^2 \frac{\Phi_M}{\Phi_H}
\]  

(3.37)

The Prandtl number \( Pr = \Phi_M/\Phi_H \) is evaluated at \( z = 0.1z_i \).

- **Entrainment zone**. Entrainment at the top of the convective boundary layer is taken into account explicitly. The buoyant flux at \( z = z_i \) is assumed to be proportional to the surface heat flux:

\[
(Q_v)_{z=z_i} = -C_{entr}Q_{0v}
\]  

(3.38)

where the entrainment constant \( C_{entr} \) is determined from experimental data. The numerical value of 0.2 is chosen from Driedonks and Tennekes (1984).

Knowing the flux at the top of the mixed layer, the exchange coefficient can be expressed as:

\[
\tilde{K}_H = -C_{entr}Q_{0v} \frac{\Delta z}{\Delta \theta_v} \quad \text{with}
\]

\[
(\Delta \theta_v)_{k+1/2} = \frac{1}{c_{dry}} \{ s_k - s_{k+1} - 0.5(\delta - \epsilon)(q_k - q_{k+1})(s_k + s_{k+1}) \}
\]  

(3.39)

Then at \( z = z_i \) :

\[
K_H = \max(K_H, \tilde{K}_H)
\]  

(3.40)
Instead of the exchange coefficients \((K)_{k+1/2}\) themselves, the scaled quantities \((K)^*_{k+1/2}\) are computed:

\[
K^*_{k+1/2} = K_{k+1/2} \alpha \rho_{k+1/2} \left[ \frac{2g\Delta t}{(\Delta z)^2_{k+1/2}} \right]
\]

where \(\alpha\) is the implicitness factor of the finite difference scheme (see equation (3.44)).

### 3.4 Solution of the Vertical Diffusion Equations

The equations for turbulent transfer are solved with the tendencies from the adiabatic (subscript ‘dyn’) and radiative processes (subscript ‘rad’) as source terms in the right hand side:

\[
\frac{\partial \psi}{\partial t} = \frac{\partial J_w}{\partial p} \bigg|_{\text{dyn}} + \frac{\partial \psi}{\partial t} \bigg|_{\text{rad}}
\]

Since the thickness of the model layers \(\Delta z\) is small near the ground, the time-stepping procedure must be implicit in order to avoid numerical instability when \(K\Delta t/(\Delta z)^2 > 1\). However, the interaction with the adiabatic and radiative processes is treated implicitly, and Janssen et al. (1992) have shown that if the tendencies are not added to the right hand side of equation (3.42) a time step dependent equilibrium, and a too low numerical drag coefficient for high wind speeds, arise. By applying a ‘fractional-steps’ method (Beljaars, 1991), the discretization of equation (3.42) becomes, for \(1 < k < l\).
\[
\frac{\partial \psi}{\partial t}_{\text{diff}} = \frac{\psi^{t+1} - \{ \psi^t + \Delta \psi_{\text{rad}} + \Delta \psi_{\text{dyn}} \}}{\Delta t}
\]

\[
= \frac{g}{p_{k+1/2} - p_{k-1/2}} \left\{ \rho_{k+1/2} K^t + \frac{K^t}{\alpha} \left( \psi_{k+1} - \psi_k \right) - \rho_{k-1/2} K^t \frac{1}{\alpha} \left( \psi_k - \psi_{k-1} \right) \right\}
\]

where

\[
\hat{\psi} = \alpha \psi^{t+1} + (1 - \alpha) \psi^t
\]

The parameter \( \alpha \) determines the implicitness of the scheme. For \( \alpha = 0 \) the scheme is explicit, for \( \alpha = 0.5 \) we have a Crank–Nicholson and for \( \alpha = 1 \) we have an implicit backward scheme. In the model, \( \alpha = 1.5 \), to avoid non-linear instability from the \( K \)-coefficients. The exchange coefficients are computed from the mean variables at \( t - 1 \).

The previous equation can be written as

\[
\frac{K^t_{k-1/2}}{\Delta p_k} \left( \frac{\psi_{k-1}}{\alpha} \right) + \left[ 1 + \frac{K^t_{k+1/2}}{\Delta p_k} + \frac{K^t_{k-1/2}}{\Delta p_k} \right] \frac{\dot{\psi}_k}{\alpha} - \frac{K^t_{k+1/2}}{\Delta p_k} \frac{\dot{\psi}_{k+1}}{\alpha} = \frac{\psi^f}{\alpha} + \Delta \psi_{\text{rad}} + \Delta \psi_{\text{dyn}}
\]

leading to the inversion of a tridiagonal matrix to solve for \( \dot{\psi}/\alpha \). The coefficients \( K^t \) are defined from (3.41).

At the lowest level (\( k = l \)) the equation includes the surface fluxes which are obtained by averaging over \( N_T \) tiles:

\[
\frac{\dot{\psi}_l}{\alpha} = \frac{\psi^f}{\alpha} + \Delta \psi_{\text{rad}} - K^t_{l-1/2} \frac{\psi_l - \psi_{l-1}}{\alpha \Delta p_l} + \sum_{i=1}^{N_T} \frac{C_{wi}^*}{\alpha \Delta p_l} \left( A_{li} \dot{\psi}_l - A_{\text{surf}i} \dot{\psi}_{\text{surf}i} \right)
\]

with \( C_{wi}^* = C_{wi}^{t-1} \left| U_i \right| g \rho \alpha \Delta t \) and

\[
\begin{align*}
\psi &= 0 & A_{li} &= 1 & A_{\text{surf}i} &= 1 & N_T &= 1 \text{ for } \psi = u, v \\
\psi_{\text{surf}} &= s_{\text{skin}i} & A_{li} &= 1 & A_{\text{surf}i} &= 1 & N_T &= 8 \text{ for } \psi = s \\
\psi_{\text{surf}} &= q_{\text{sut}}(T_{\text{skin}i}) & A_{li} &= \alpha_{li} & A_{\text{surf}i} &= \alpha_{\text{surf}i} & N_T &= 8 \text{ for } \psi = q
\end{align*}
\]

Eq. (3.46) can be re-written

\[
\frac{K^t_{l-1/2}}{\Delta p_l} \left( \frac{\psi_{l-1}}{\alpha} \right) + \left[ 1 + \sum_{i=1}^{N_T} F_i C_{wi}^* A_{li} \frac{\psi_l}{\alpha} + \frac{K^t_{l-1/2}}{\Delta p_l} \frac{\psi_l}{\alpha} \right] = \frac{\psi^f}{\alpha} + \Delta \psi_{\text{rad}} - \sum_{i=1}^{N_T} F_i C_{wi}^* A_{\text{surf}i} \frac{\psi_{\text{surf}i}}{\alpha}
\]
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The implicit value at the surface \( \psi_{\text{surf}} \) on the right hand side is obtained from coupling this last equation with the calculation of the surface energy budget through the computation of the skin surface temperature (see equation (3.62) in Section 3.5).

At the top of the atmosphere \((k = 1)\) turbulent fluxes at set to zero and we have

\[
\frac{\dot{\psi}_1}{\alpha} = \frac{\dot{\psi}_1}{\alpha} + \Delta \psi_{\text{dy}} + \Delta \psi_{\text{rad}} - K^{*}_{-1/2} \frac{\psi_1 - \psi_0}{\alpha \Delta p_1} \tag{3.49}
\]

which can be re-written

\[
-\frac{K^{*}_{1/2} \dot{\psi}_0}{\Delta p_1 \alpha} + \left( 1 + \frac{K^{*}_{1/2}}{\Delta p_1} \right) \frac{\dot{\psi}_1}{\alpha} = \frac{\dot{\psi}_1}{\alpha} + \Delta \psi_{\text{dy}} + \Delta \psi_{\text{rad}} \tag{3.50}
\]

The tridiagonal matrix equation is solved by a downward elimination scan followed by an upward back substitution (Press et al., 1992, pp 42-43).

3.5 The skin temperature

The surface energy balance is satisfied independently for the tiles by calculating the skin temperature for each tile. The skin layer represents the vegetation layer, the top layer of the bare soil, or the top layer of the snow pack, has no heat capacity and therefore responds instantaneously to changes in e.g. radiative forcing. In order to calculate the skin temperature, the surface energy balance equation is linearized for each tile leading to an expression for the skin temperature. This procedure is equivalent to the Penmann-Monteith approach which can be derived by eliminating the skin temperature from the surface energy balance equation assuming that the net radiation minus ground heat flux is known (e.g. Brutsaert, 1982). The approach followed here is an extension to the Penmann-Monteith derivation in the sense that it allows for coupling with the underlying soil (or snow, ice). Because of the short time scale associated with the skin layer, the equation for its temperature is solved implicitly together with the vertical diffusion in the boundary layer.

The following general discussion applies to each tile but the parameters are tile dependent as discussed in the land surface part of the documentation (Chapter 7). The surface energy balance equation can be written as:

\[
\mathcal{R}_{SW} + \mathcal{R}_{LW} + H + L J_q = \Lambda_{\text{skin}} (T_{\text{skin}} - T_s) \tag{3.51}
\]

where \( \mathcal{R}_{SW} \) and \( \mathcal{R}_{LW} \) are the net shortwave and longwave radiation fluxes at the surface and the right hand side represents the ground heat flux through coupling with the underlying soil, snow or ice with temperature \( T_s \). The turbulent sensible and latent heat fluxes are

\[
H = J_s - c_{\text{dry}} T_{\text{skin}} \delta J_q \tag{3.52}
\]

\[
J_s = \rho C_h [U_s] \{ s_l - s_{\text{skin}} \} \tag{3.53}
\]

\[
J_q = \rho C_q [U_s] \{ \alpha_l q_l - \alpha_{\text{surf}} q_{\text{sat}}(T_{\text{skin}}) \} \tag{3.54}
\]

In order to solve for the skin temperature implicitly, the surface energy balance is solved together with the vertical diffusion equations. After the downward elimination scan of the tridiagonal system of equations (3.45) a relation
is obtained between the lowest model level values and the surface values, i.e. between $\dot{q}_t$ and $\dot{q}_{sat}(\dot{T}_{skin})$, and between $\dot{s}_t$ and $\dot{s}_{skin}$:

$$\dot{s}_t = Z_{AS}\dot{s}_{skin} + Z_{BS}$$  \hspace{1cm} (3.55)

$$\dot{q}_t = Z_{AQ}\dot{q}_{sat}(\dot{T}_{skin}) + Z_{BQ}$$  \hspace{1cm} (3.56)

Since the vertical diffusion equation is formulated in terms of the time extrapolated parameters (indicated by a hat, see equation (3.44)), the skin temperature has to be extrapolated as well. Eliminating the lowest model level parameters and linearizing with respect to previous time step skin temperature leads to

$$H = \rho C_H[U]l\{(Z_{AS} - 1)\dot{s}_{skin} + Z_{BS}\} - c_{pdry}\dot{T}_{skin}\delta J_q$$  \hspace{1cm} (3.57)

$$J_q = \rho C_Q[U]l\left\{a_I(Z_{AQ} - a_{surf})\left[q_{sat}(T^d_{skin}) + \frac{\partial q_{sat}}{\partial T}(\dot{T}_{skin} - T^d_{skin})\right] + a_IZ_{BQ}\right\}$$  \hspace{1cm} (3.58)

Also $\dot{s}_{skin}$ needs to be expressed in surface variables. For this purpose the moisture correction in $c_p$ is evaluated from the previous time level:

$$\dot{s}_{skin} = c_p\dot{T}_{skin}$$ \hspace{1cm} (3.59)

$$c_p = c_{pdry}\left\{1 + \delta a_{surf}\dot{q}_{sat}(T^d_{skin}) + \delta(1 - \alpha_I)q^d_I\right\}$$ \hspace{1cm} (3.60)

The net long-wave radiation at the surface is linearized with respect to skin temperature at the previous radiation time step (indicated by superscript $\text{trad}$, which can be up to 3 hours earlier):

$$\dot{R}_{LW} = \dot{R}_{LW}^{\text{trad}} - 4(T^d_{skin})^3(\dot{T}_{skin} - T^{\text{trad}}_{skin})$$ \hspace{1cm} (3.61)

Substituting $\dot{s}_{skin}$ in (3.57) and replacing $H$ and $J_q$ in surface energy balance equation (3.51) by equations (3.57) and (3.58) leads to an expression for skin temperature $\dot{T}_{skin}$ at the extrapolated time level.

$$\dot{T}_{skin} = \left[\Lambda_{skin}T_s + \dot{R}_{SW} + \dot{R}_{LW}^{\text{trad}} + 4(T^{\text{trad}}_{skin})^4 + \rho C_H[U]Z_{BS} + \rho C_Q(L - c_{pd}\dot{T}_{skin}^d)(U)\left\{(a_IZ_{AQ} - a_{surf})\left[q_{sat}(T^d_{skin}) - \frac{\partial q_{sat}}{\partial T}(\dot{T}_{skin} - T^d_{skin})\right] + a_IZ_{BQ}\right\} \right]^{-1}$$ \hspace{1cm} (3.62)

with $c_p$ from equation (3.60). Following the downward elimination scan of the tridiagonal matrices for $s$ and $q$, equation (3.62) is solved for all the tiles, using the appropriate parameters for each tile (note that also the transfer coefficients and therefore coefficients $Z_{AQ}$, $Z_{BQ}$, $Z_{AS}$, $Z_{BS}$ are tile dependent). The resulting skin temperatures are used in (3.48) with the corresponding weights of the tiles as a boundary condition before doing the upward scanning back-substitution.
This procedure is fully implicit for the dominant tile in the sense that atmospheric and skin variables are in equilibrium at the new time level. However, equilibrium for non-dominant tiles is not necessarily achieved. It can happen that the surface fluxes from the dominant tile changes the temperature and moisture substantially at the lowest model level. If the fluxes to another tile (with small fraction) happen to be very different, this tile will not see the correct atmospheric state in the computation of the skin temperature. A full implicit coupling would require the solution of a matrix problem involving the skin temperatures of all the tiles simultaneously.

### 3.6 TENDENCY CALCULATIONS AND ENERGY DISSIPATION

Total wind and specific humidity tendencies after the vertical diffusion (including also the dynamics and the radiation) are

\[
\begin{align*}
\frac{\partial u}{\partial t} &= \frac{u^{t+1} - u^{t}}{\Delta t} \\
\frac{\partial v}{\partial t} &= \frac{v^{t+1} - v^{t}}{\Delta t} \\
\frac{\partial q}{\partial t} &= \frac{q^{t+1} - q^{t}}{\Delta t}
\end{align*}
\]  

(3.63)

The kinetic energy lost by the mean flow through the diffusion process, \(E_{\text{diff}}\), is

\[
E_{\text{diff}} = 2\Delta t \cdot \left[ \frac{u^{t+1} + u^{t}}{2} \right] + 2\Delta t \cdot \left[ \frac{v^{t+1} + v^{t}}{2} \right]
\]  

(3.64)

The kinetic energy lost is assumed to be transformed locally into internal energy. This procedure by-passes the sub-grid scale energy cascade, but it allows to have a closed energy cycle in the model (the term is generally small)

\[
\frac{\partial T}{\partial t} \bigg|_{\text{diff + dyn + rad}} = \frac{1}{2M} \left[ \frac{s^{t+1} - \phi + E_{\text{diff}}}{C_{p, dry}(1 + eq^{t+1})} - T^{t+1} \right]
\]  

(3.65)

### 3.7 SHORTER TIME STEP IN THE VERTICAL DIFFUSION SCHEME

The vertical diffusion scheme is called three times in every physics time step, with a time step of 1/3 of the standard time step.

### 3.8 DIAGNOSTIC COMPUTATIONS FOR POSTPROCESSING

#### 3.8.1 Diagnostic boundary layer height

Because of its importance for applications (e.g. in air pollution modelling), the boundary layer height is diagnosed and made available for postprocessing. The parametrization of the mixed layer (and entrainment) already uses a model level index as boundary layer height, but in order to get a continuous field, also in neutral and stable situations the parcel lifting method (or bulk Richardson method) proposed by Troen and Mahrt (1986) is used as a diagnostic, independent of the turbulence parametrization. Boundary layer height \(h_{BL}\) is defined as the level where
the bulk Richardson number, based on the difference between quantities at that level and the lowest model level, reaches the critical value $Ri_{cr} = 0.25$. The bulk Richardson is computed from the following set of equations

\[
|\Delta U|^2 = (u_{hbl} - u_I)^2 + (v_{hbl} - v_I)^2
\]

\[
s_{sl} = c_{pdr}T_{sl}(1 + eq_I + gz_l)
\]

\[
s_{vhbl} = c_{pdr}T_{hbl}(1 + eq_{hbl}) + gh_{hbl}
\]

\[
\Delta s = 8.5c_{pdr}u_*Q_{sbl}/w_s
\]

\[
w_s = \left\{ u_*^3 + 0.6g/TQ_{sbl}h_{hbl}\right\}^{1/3} \quad \text{unstable}
\]

\[
w_s = u_* \quad \text{stable}
\]

\[
R_i_b = h_{hbl}\frac{2g(s_{vhbl} - s_{sl} - \Delta s)}{(s_{vhbl} + s_{sl} - gh_{hbl} - gz_l)|\Delta U|^2}
\]

where index $l$ indicates the lowest model level and $h_{hbl}$ indicates the boundary layer height i.e the level where $Ri_b = Ri_{cr}$. The virtual dry static energy from the lowest level $s_{sl}$ is increased with a turbulent part $\Delta s$ and compared to the virtual dry static energy at boundary layer height $h_{hbl}$. The boundary layer height is found by a vertical scan from the surface upwards. If the boundary layer height is found to be between two levels a linear interpolation is done to find the exact position. Since the boundary layer height is needed for $w_s$, the upward scan is done twice. The first one uses $h_{BL} = 1000m$ in the expression for $w_s$; the second scan uses the result of the first scan.

### 3.8.2 Wind at 10 m level

Wind at the 10 m level is computed for postprocessing because it is the standard level for SYNOP observations. It can be obtained rather easily by vertical interpolation between the lowest model level and the surface, making use of profile functions (3.7) and (3.8). This procedure is appropriate over the ocean or in areas where the surface is smooth and homogeneous. However, the postprocessed field is meant to be comparable to wind from SYNOP observations and for observations over land WMO requires SYNOP stations to be in open terrain in order to be well exposed to wind. So the SYNOP wind observations are not necessarily compatible with the wind that is representative for a large area (i.e. a grid box from the model). Over inhomogeneous terrain, the problem can be particularly serious, because the “aerodynamic roughness length” in the model is adjusted to provide sufficient drag at the surface which is dominated by the rough elements. This approach leads to a low area-averaged wind speed which is not comparable to the “open-terrain” wind speed as observed by WMO stations.

In order to make the postprocessed wind compatible with SYNOP observations, the concept of exposure correction is introduced. The open-terrain wind is obtained by taking the wind information from such a height above the surface that it is less influenced by the underlying terrain. This height is called the blending height $h_{blend}$, and for the the interpolation to 10 m an aerodynamic roughness length is used that is typical for open terrain with grassland.

The interpolation procedure is as follows. First the blending height and the interpolation roughness length are set dependent on the model roughness length field:

\[
h_{blend} = 75, \quad z_{0MWMO} = 0.03, \quad F_{blend} = (u_{blend}^2 + v_{blend}^2)^{1/2} \quad \text{if } z_{0M} > 0.03
\]

\[
h_{blend} = z_l, \quad z_{0MWMO} = z_{0M}, \quad F_{blend} = (u_l^2 + v_l^2)^{1/2} \quad \text{if } z_{0M} < 0.03
\]

\[
(3.67)
\]
\[ F_{10} = F_{\text{blend}} \frac{\log\left( \frac{z_1 + z_{0\text{WMWO}}}{z_{0\text{WMWO}}} \right) - \Psi_M\left( \frac{z_1 + z_{0\text{WMWO}}}{L} \right) + \Psi_M\left( \frac{z_{0\text{WMWO}}}{L} \right)}{\log\left( \frac{z_{\text{blend}} + z_{0\text{WMWO}}}{z_{0\text{WMWO}}} \right) - \Psi_M\left( \frac{z_{\text{blend}} + z_{0\text{WMWO}}}{L} \right) + \Psi_M\left( \frac{z_{0\text{WMWO}}}{L} \right)} \]  

(3.68)

where \( z_{10} = 10 \text{ m} \), \( F_{\text{blend}} \) is the horizontal wind speed at the blending height either interpolated from model levels to 75 m or copied from the lowest model level, and \( F_{10} \) is the resulting horizontal windspeed at 10 m. The wind speed from equation (3.68) is converted to components making use of the wind direction from the lowest model level.

### 3.8.3 Temperature and humidity at the 2 m level

Computation of temperature and moisture at the 2 m level is based on interpolation between the lowest model level and the surface making use of the same profile functions as in the parametrization of the surface fluxes. The following expressions are derived from equations (3.9) and (3.10)

\[ s_2 = s_{\text{surf}} + (s_1 - s_{\text{surf}}) \frac{\log\left( \frac{z_2 + z_{0\text{WMWO}}}{z_{0\text{WMWO}}} \right) - \Psi_H\left( \frac{z_2 + z_{0\text{WMWO}}}{L} \right) + \Psi_H\left( \frac{z_{0\text{WMWO}}}{L} \right)}{\log\left( \frac{z_1 + z_{0\text{WMWO}}}{z_{0\text{WMWO}}} \right) - \Psi_H\left( \frac{z_1 + z_{0\text{WMWO}}}{L} \right) + \Psi_H\left( \frac{z_{0\text{WMWO}}}{L} \right)} \]  

(3.69)

\[ q_2 = q_{\text{surf}} + (q_1 - q_{\text{surf}}) \frac{\log\left( \frac{z_2 + z_{0\text{WMWO}}}{z_{0\text{WMWO}}} \right) - \Psi_H\left( \frac{z_2 + z_{0\text{WMWO}}}{L} \right) + \Psi_H\left( \frac{z_{0\text{WMWO}}}{L} \right)}{\log\left( \frac{z_1 + z_{0\text{WMWO}}}{z_{0\text{WMWO}}} \right) - \Psi_H\left( \frac{z_1 + z_{0\text{WMWO}}}{L} \right) + \Psi_H\left( \frac{z_{0\text{WMWO}}}{L} \right)} \]  

(3.70)

with \( z_2 = 2 \text{ m} \), \( z_{0\text{WMWO}} = z_{0\text{QWMWO}} = 0.003 \) if \( z_{0M} > 0.03 \), and otherwise \( z_{0\text{WMWO}} = z_{0H} \) and \( z_{0\text{QWMWO}} = z_{0Q} \). Temperature \( T_2 \) is derived from \( s_2 \) and \( q_2 \) with equation (3.3). Also the dew point is computed from \( q_2 \) and surface pressure. The dew point uses the saturation formulation with respect to water to be consistent with WMO reporting practise. If the resulting dew point is lower than temperature \( T_2 \), the dew point is set equal to temperature.

### 3.8.4 Wind gusts

The computation of gusts is intended to be compatible with WMO observing practise for wind extremes. In order to get uniform observations, WMO defines a wind gust as the maximum of the wind averaged over 3 second intervals.

First the friction and the horizontal wind speed at the 10 m level are computed from the lowest model level (no exposure correction)

\[ u_* = \frac{F_l}{\frac{\log\left( \frac{z_1 + z_{0M}}{z_{0M}} \right) - \Psi_M\left( \frac{z_1 + z_{0M}}{L} \right) + \Psi_M\left( \frac{z_{0M}}{L} \right)}{\log\left( \frac{z_{10} + z_{0M}}{z_{0M}} \right) - \Psi_M\left( \frac{z_{10} + z_{0M}}{L} \right) + \Psi_M\left( \frac{z_{0M}}{L} \right)}} \]  

(3.71)

\[ F_{M10} = \frac{u_*}{\kappa} \left[ \frac{\log\left( \frac{z_{10} + z_{0M}}{z_{0M}} \right) - \Psi_M\left( \frac{z_{10} + z_{0M}}{L} \right) + \Psi_M\left( \frac{z_{0M}}{L} \right)}{\log\left( \frac{z_{10} + z_{0M}}{z_{0M}} \right) - \Psi_M\left( \frac{z_{10} + z_{0M}}{L} \right) + \Psi_M\left( \frac{z_{0M}}{L} \right)} \right] \]
To simulate gusts, the standard deviation of the horizontal wind is estimated on the basis of the similarity relation by Panofsky et al. (1977)

\[
\sigma_u = 2.29u_* \left( \frac{1 + 0.5z_i}{L} \right)^{1/3} \quad \text{for } L < 0
\]
\[
\sigma_u = 2.29u_* \quad \text{for } L > 0
\]  \hspace{1cm} (3.72)

with \( z_i = 1000m \). The difference between the gust and \( F_{\text{M10}} \) is proportional to \( \sigma_u \), where the multiplier has been determined from universal turbulence spectra for a 50% exceeding probability of the three-second wind gust (see Beljaars, 1987). The resulting wind gust is

\[
F_{\text{gust}} = F_{\text{M10}} + C_{\text{ugn}} \sigma_u
\]  \hspace{1cm} (3.73)

with parameter \( C_{\text{ugn1}} = C_{\text{ugn}} \) for open terrain and \( C_{\text{ugn}} = 7.71 \). To correct for extremem gusts over mountainous terrain, where the roughness length is extreme, the following expression is used

\[
C_{\text{ugn1}} = \frac{1}{\kappa} \ln \left( \frac{5 \exp(\kappa C_{\text{ugn}}) + z_{0M}}{5 + z_{0M}} \right)
\]

From the controlling parameters it is clear that the effects of surface friction (through surface roughness) and stability are captured. However, the approach is not adequate for gusts in baroclinic situations and gusts due to strong convective events. Parameter \( F_{\text{gust}} \) is computed every time step and its maximum since the last postprocessing time is written out for archiving.

### 3.9 Code

Vertical diffusion, which affects temperature, velocities and specific humidity, is performed in subroutine VDFMAIN called by VDFOUTER which, in turn, is called by CALLPAR. VDFOUTER calls VDFMAIN three times with \( 1/3 \) of the normal time step (these two routines have identical arguments so VDFMAIN can be called directly if the vertical diffusion is only needed once per time step).

At the start of the model integration the following setup routines are called to initialize modules specific to the vertical diffusion code:

- **SUVDF.** Setup routine for a number of parametrization constants.
- **SUVDFS.** Setup routine for constants and tables related to the stability functions. Stability functions are included as statement functions from fcvds.h.
- **SUVEG.** Setup routine for vegetation and tile parameters.

The main subroutine (VDFMAIN) does a sequence of computations and subroutine calls:

- The tiled surface fluxes and tiled skin temperatures are cycled from time step to time step (fluxes are needed for the first guess of stability parameters), but are not available at the start of the forecast. For the first time step, neutral transfer coefficients are used to estimate momentum fluxes, the tiled skin temperatures are set equal to the grid box averaged skin temperature from the initial condition, and the sensible and latent heat fluxes needed as a first guess for the Obukhov length computation are set to zero.
Chapter 3 ‘Turbulent diffusion and interactions with the surface’

- **VFDUPDZ0.** This routine computes roughness lengths for momentum, heat and moisture over ocean surfaces according to equation (3.24). It also computes surface buoyancy flux and Obukhov length from the fluxes of the previous time level.
- The dry static energy is computed on model levels.
- A grid box average of the surface albedo is computed from the tile albedo and the tile fractions.
- **VDFSURF.** This routine prepares the surface boundary conditions for temperature and humidity and is called for every tile. The following quantities are computed: the surface specific humidity at saturation, the derivative of the saturation humidity curve at the surface, surface dry static energy, and vegetation stomatal resistances (see Chapter 7).
- **VFDSEXCS.** This routine determines the drag transfer coefficients between the surface and the lowest model level with the thermal stability expressed as function of the Obukhov length. It is called for every tile. The implicit relation between \(z/L\) and the Richardson number \(R_i_{\text{bulk}}\) is solved iteratively (using the Newton method with the derivative approximated in finite differences). Pre-computed tables defined in subroutine **SUVDFF** are used to obtain the first guess in stable conditions (\(R_i>0\)) at the first time step. Transfer coefficients are multiplied by a constant factor \(\rho a_2\Delta tg\).
- **VDFEVP.** This routine computes for each tile the equivalent evapo-transpiration efficiency and the corresponding parameters \(a_t\) and \(a_{\text{surf}}\) defined by the land surface scheme (see chapter 7). Dry static energy at the surface at time level \(t\) is estimated as well.
- **VDFSFLX.** This routine computes surface fluxes for each tile (heat flux, evaporation, momentum flux and virtual heat flux) at time \(t\) for later use in similarity functions and for the estimation of the diagnostic boundary layer depth.
- **VDFDPBL.** This routine diagnoses the boundary layer height for time level \(t\). This boundary layer height is for postprocessing only and is not used by the parametrization.
- **VFDSEXCU.** This routine determines the turbulent diffusion coefficients between the model levels above the surface layer. In unstable surface conditions, the depth of a mixed layer is estimated where the diffusion coefficients are expressed according to equations (3.36) and (3.37). Above the mixed layer, the diffusion coefficients are expressed from local similarity theory with equations (3.31) if \(R_i<0\). In layers with \(R_i>0\), diffusion coefficients are expressed as function of the Richardson number according to equation (3.32). The entrainment rate at the top of the mixed layer is computed according to (3.40).
- **VDFDFIM.** This routine solves the diffusion equation for momentum, by Gaussian elimination of the tridiagonal matrices.
- **VDFDFIF.** This routine solves the diffusion equations for dry static energy and specific humidity. A downward elimination scan is done through the tridiagonal matrices, and coefficients \(z_{AS}, z_{BS}, z_{AQ}\), and \(z_{BQ}\) are computed for each tile. Then, subroutine **VDFTSK** is called for each tile to compute the skin temperatures from the surface energy balance equation. Subsequently the tiled skin temperatures are used as a boundary condition and the back-sustitution is performed.
- **VDFINCR.** This routine computes the tendencies of the prognostic variables and estimates the kinetic energy dissipation.
- **VDFDFLX.** This routine computes the tile fluxes at the new time levels. These are also the fluxes to be used in the land surface scheme. Averaging over the tiles is also done for diagnostic purposes and postprocessing.
- **VDFPPCPL.** This routine computes the surface 2 metre temperature and humidity (dew point and specific humidity), and the wind at 10 m.
- **VDFPPGUST.** This routine computes wind gusts as they are typically observed by standard WMO SYNOP stations.
APPENDIX A LIST OF SYMBOLS

\( C_H \)  transfer coefficient for heat
\( C_M \)  transfer coefficient for momentum (drag coefficient)
\( C_Q \)  transfer coefficient for moisture
\( c_p \)  specific heat at constant pressure of moist air
\( c_{pdry} \)  specific heat at constant pressure of dry air
\( c_{pvap} \)  specific heat at constant pressure of water vapour
\( E_{diss} \)  kinetic energy lost by the diffusion process
\( f \)  Coriolis parameter
\( F_{blend} \)  horizontal wind speed at blending height (for pp of 10 m wind)
\( F_{10} \)  horizontal wind speed at 10 m level (for pp)
\( g \)  acceleration of gravity
\( h_{HL} \)  diagnosed boundary layer height
\( h_{blend} \)  blending height (for pp of 10 m wind)
\( J_q \)  vertical turbulent flux of \( \psi \)
\( J_q \)  surface humidity flux
\( J_x \)  surface flux of dry static energy
\( J_M \)  surface momentum flux
\( K_q \)  turbulent exchange coefficient for \( \psi \)
\( K_H \)  turbulent exchange coefficient for heat
\( K_M \)  turbulent exchange coefficient for momentum
\( K_Q \)  turbulent exchange coefficient for moisture
\( L \)  Obukhov length
\( L \)  latent heat of vaporization/sublimation
\( l_H \)  mixing length for heat
\( l_M \)  mixing length for momentum
\( N_T \)  number of tiles
\( p \)  pressure
\( Pr \)  Prandtl number
\( q \)  specific humidity
\( q^* \)  \( = J_q/(\rho u^*) \)
\( Q_{lv} \)  virtual temperature flux in the surface layer
\( q_{sat} \)  saturation specific humidity
\( R_{dry} \)  gas constant for dry air
\( R_{vap} \)  gas constant for water vapour
\( R_{LW} \)  net long wave radiation at the surface
\( R_{SW} \)  net short wave radiation at the surface
\( RH_{surf} \)  relative humidity at the surface
\( Ri \)  local Richardson number
\( Ri_{bulk} \)  bulk Richardson number for the surface layer
\( s \)  dry static energy
\( s_c \)  virtual dry static energy
\( s^* \)  \( = J_q/(\rho u^*) \)
\( T \)  temperature
\( t \)  time

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**Part IV: ‘Physical processes’**

**APPENDIX A LIST OF SYMBOLS**

- \( C_H \): transfer coefficient for heat
- \( C_M \): transfer coefficient for momentum (drag coefficient)
- \( C_Q \): transfer coefficient for moisture
- \( c_p \): specific heat at constant pressure of moist air
- \( c_{pdry} \): specific heat at constant pressure of dry air
- \( c_{pvap} \): specific heat at constant pressure of water vapour
- \( E_{diss} \): kinetic energy lost by the diffusion process
- \( f \): Coriolis parameter
- \( F_{blend} \): horizontal wind speed at blending height (for pp of 10 m wind)
- \( F_{10} \): horizontal wind speed at 10 m level (for pp)
- \( g \): acceleration of gravity
- \( h_{HL} \): diagnosed boundary layer height
- \( h_{blend} \): blending height (for pp of 10 m wind)
- \( J_q \): vertical turbulent flux of \( \psi \)
- \( J_q \): surface humidity flux
- \( J_x \): surface flux of dry static energy
- \( J_M \): surface momentum flux
- \( K_q \): turbulent exchange coefficient for \( \psi \)
- \( K_H \): turbulent exchange coefficient for heat
- \( K_M \): turbulent exchange coefficient for momentum
- \( K_Q \): turbulent exchange coefficient for moisture
- \( L \): Obukhov length
- \( L \): latent heat of vaporization/sublimation
- \( l_H \): mixing length for heat
- \( l_M \): mixing length for momentum
- \( N_T \): number of tiles
- \( p \): pressure
- \( Pr \): Prandtl number
- \( q \): specific humidity
- \( q^* \): \( = J_q/(\rho u^*) \)
- \( Q_{lv} \): virtual temperature flux in the surface layer
- \( q_{sat} \): saturation specific humidity
- \( R_{dry} \): gas constant for dry air
- \( R_{vap} \): gas constant for water vapour
- \( R_{LW} \): net long wave radiation at the surface
- \( R_{SW} \): net short wave radiation at the surface
- \( RH_{surf} \): relative humidity at the surface
- \( Ri \): local Richardson number
- \( Ri_{bulk} \): bulk Richardson number for the surface layer
- \( s \): dry static energy
- \( s_c \): virtual dry static energy
- \( s^* \): \( = J_q/(\rho u^*) \)
- \( T \): temperature
- \( t \): time
Chapter 3 ‘Turbulent diffusion and interactions with the surface’

\[ |U| \] horizontal wind speed
\[ u, v \] horizontal wind components
\[ u_* \] friction velocity = \( (J_M/\rho)^{1/2} \)
\[ w_* \] free convection velocity scale
\[ w_{turb} \] turbulent velocity scale
\[ z_{0M} \] roughness length for momentum (aerodynamic roughness length)
\[ z_{0H} \] roughness length for heat
\[ z_{0Q} \] roughness length for moisture
\[ z_i \] scale height of the boundary layer
\[ z_l \] height of the lowest model level \( l \)
\[ z_{0M\text{MWMO}} \] roughness length for momentum at SYNOP station
\[ z_{0H\text{HWMO}} \] roughness length for heat at SYNOP station
\[ z_{0Q\text{QWMO}} \] roughness length for moisture at SYNOP station
\[ z_2 \] height of screen level observation (2 m)
\[ z_{10} \] height of surface wind observation (10 m)
\[ \alpha \] implicitness factor for diffusion equation
\[ \alpha_{Ch} \] Charnock parameter
\[ \beta \] scaling parameter for asymptotic mixing length
\[ \Delta t \] time step
\[ \Delta z \] vertical grid length
\[ \delta \] \( e_{\text{pewp}}/e_{\text{pdry}} - 1 \)
\[ \varepsilon \] \( (R_{\text{wpc}}/R_{\text{dry}}) - 1 \)
\[ \theta_v \] virtual potential temperature
\[ \kappa \] Von Kármán’s constant
\[ \lambda \] asymptotic mixing length
\[ \Lambda_{\text{skin}} \] conductivity of
\[ \nu \] kinematic viscosity
\[ \rho \] density
\[ \sigma_u \] standard deviation of horizontal wind
\[ \zeta = z/L \] geopotential
\[ \phi \] symbolic reference to a conservative quantity
\[ \Phi_M \] universal gradient stability function for wind
\[ \Phi_H \] universal gradient stability function for temperature
\[ \Phi_Q \] universal gradient stability function for moisture
\[ \psi \] symbolic reference to a conservative quantity
\[ \Psi_M \] universal profile stability function for wind
\[ \Psi_H \] universal profile stability function for temperature
\[ \Psi_Q \] universal profile stability function for moisture

Subscripts:
\[ i \] tile index
\[ k \] level index (counted from model top downwards)
\[ l \] referring to lowest model level
\[ \text{skin} \] referring to the skin layer
\[ \text{surf} \] referring to the surface
Superscripts:
- $t$: index for old time level, indicating beginning of time step
- $t+1$: index for new time level, indicating end of time step
- $\text{trad}$: index referring to the latest full radiation time step

Special symbols:
- $\psi$: implicit variable $\psi$ defined by equation (3.44)
Part IV: PHYSICAL PROCESSES

CHAPTER 4 Subgrid-scale orographic drag

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4.1 General principles
4.2 Description of the scheme
  4.2.1 Blocked-flow drag
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  4.4.1 GWSETUP
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4.1 GENERAL PRINCIPLES

The influence of subgrid-scale orography on the momentum of the atmosphere, and hence on other parts of the physics, is represented by a combination of lower-troposphere drag created by orography assumed to intersect model levels, and vertical profiles of drag due to the absorption and/or reflection of vertically propagating gravity waves generated by stably stratified flow over the subgrid-scale orography. The scheme is described in detail in Lott and Miller (1996).

The scheme is based on ideas presented by Baines and Palmer (1990), combined with ideas from bluff-body dynamics. The assumption is that the mesoscale-flow dynamics can be described by two conceptual models, whose relevance depends on the non-dimensional height of the mountain via.

\[ H_n = \frac{NH}{U} \]  

(4.1)

where \( H \) is the maximum height of the obstacle, \( U \) is the wind speed and \( N \) is the Brunt–Väisälä frequency of the incident flow.

At small \( H_n \) all the flow goes over the mountain and gravity waves are forced by the vertical motion of the fluid. Suppose that the mountain has an elliptical shape and a height variation determined by a parameter \( b \) in the along-ridge direction and by a parameter \( a \) in the cross-ridge direction, such that \( \gamma = a/b \leq 1 \), then the geometry of the mountain can be written in the form

\[ h(x, y) = \frac{H}{1 + x^2/a^2 + y^2/b^2} . \]

(4.2)

In the simple case when the incident flow is at right angles to the ridge the surface stress due to the gravity wave
has the magnitude

\[ \tau_{\text{wave}} = \rho_0 b G B(\gamma) NUH^2 \]

(4.3)

provided that the Boussinesq and hydrostatic approximations apply. In Eq. (4.3) \( G \) is a function of the mountain sharpness (Phillips 1984), and for the mountain given by Eq. (4.2), \( G = 1.23 \). The term \( B(\gamma) \) is a function of the mountain anisotropy, \( \gamma \), and can vary from \( B(0) = 1 \) for a two-dimensional ridge to \( B(1) = \pi/4 \) for a circular mountain.

At large \( H_n \), the vertical motion of the fluid is limited and part of the low-level flow goes around the mountain. As is explained in Section 4.2, the depth, \( Z_{\text{blk}} \), of this blocked layer, when \( U \) and \( N \) are independent of height, can be expressed as

\[ Z_{\text{blk}} = H \times \max \left( 0, \frac{H_n - H_n^{\text{crit}}}{H_n} \right) \]

(4.4)

where \( H_n^{\text{crit}} \) is a critical non-dimensional mountain height of order unity. The depth \( Z_{\text{blk}} \) can be viewed as the upstream elevation of the isentropic surface that is raised exactly to the mountain top. In each layer below \( Z_{\text{blk}} \), the flow streamlines divide around the obstacle, and it is supposed that flow separation occurs on the obstacle’s flanks.

Then, the drag, \( D_{\text{blk}}(z) \), exerted by the obstacle on the flow at these levels can be written as

\[ D_{\text{blk}}(z) = -\rho_0 C_d l(z) \frac{U|U|}{2} \]

(4.5)

Here \( l(z) \) represents the horizontal width of the obstacle as seen by the flow at an upstream height \( z \) and \( C_d \), according to the free streamline theory of jets in ideal fluids, is a constant having a value close to unity (Kirchoff 1876; Gurvitch 1965). According to observations, \( C_d \) can be nearer 2 in value when suction effects occur in the rear of the obstacle (Batchelor 1967). In the proposed parametrization scheme this drag is applied to the flow, level by level, and will be referred to as the drag of the ‘blocked’ flow, \( D_{\text{blk}} \). Unlike the gravity-wave-drag scheme, the total stress exerted by the mountain on the ‘blocked’ flow does not need to be known \textit{a priori}. For an elliptical mountain, the width of the obstacle, as seen by the flow at a given altitude \( z < Z_{\text{blk}} \), is given by

\[ l(z) = 2b \left( \frac{Z_{\text{blk}} - z}{z} \right)^{1/4} \]

(4.6)

In Eq. (4.6), it is assumed that the level \( Z_{\text{blk}} \) is raised up to the mountain top, with each layer below \( Z_{\text{blk}} \) raised by a factor \( H/Z_{\text{blk}} \). This leads, effectively, to a reduction of the obstacle width, as seen by the flow when compared with the case in which the flow does not experience vertical motion as it approaches the mountain. Then applying Eq. (4.5) to the fluid layers below \( Z_{\text{blk}} \), the stress due to the blocked-flow drag is obtained by integrating from \( z = 0 \) to \( z = Z_{\text{blk}} \), viz.

\[ \tau_{\text{blk}} = C_d \pi b \rho_0 Z_{\text{blk}} \frac{U|U|}{2} \]

(4.7)

However, when the non-dimensional height is close to unity, the presence of a wake is generally associated with upstream blocking and with a downstream foehn. This means that the isentropic surfaces are raised on the windward side and become close to the ground on the leeward side. It we assume that the lowest isentropic surface passing over the mountain can be viewed as a lower rigid boundary for the flow passing over the mountain, then the
distortion of this surface will be seen as a source of gravity waves and, since this distortion is of the same order of 
magnitude as the mountain height, it is reasonable to suppose that the wave stress will be given by Eq. (4.3), where-
ever the depth of the blocked flow, \( Z_{\text{blk}} \), although it is clearly an upper limit to use the total height, \( H \). Then, the 
total stress is the sum of a wave stress, \( \tau_{\text{wave}} \), and a blocked-flow stress whenever the non-dimensional mountain 
height \( H_n > H_{n\text{blk}} \), i.e.

\[
\tau \approx \tau_{\text{wave}} \left[ 1 + \frac{\pi C_d}{2GB(\gamma)} \max\left(0, \frac{H_n - H_{n\text{blk}}}{H_n^2}\right) \right]. \tag{4.8}
\]

The addition of low-level drag below the depth of the blocked flow, \( Z_{\text{blk}} \), enhances the gravity-wave stress term 
in Eq. (4.8) substantially.

In the present scheme the value of \( C_d \) is allowed to vary with the aspect ratio of the obstacle, as in the case of 
separated flows around immersed bodies (Landweber 1961), while at the same time setting the critical number 
\( H_{n\text{crit}} \) equal to 0.5 as a constant intermediate value. Note also that for large \( H_n \), Eq. (4.8) overestimates the drag 
in the three-dimensional case, because the flow dynamics become more an more horizontal, and the incidence of 
gravity waves is diminished accordingly. In the scheme a reduction of this kind in the mountain-wave stress could 
have been introduced by replacing the mountain height given in Eq. (4.3) with a lower ‘cut-off’ mountain height, 
\( H(H_{n\text{crit}}/H_n) \). Nevertheless, this has not been done partly because a large non-dimensional mountain height often 
corresponds to the slow flows for which the drag given by Eq. (4.8) is then, in any case, very small.

### 4.2 Description of the Scheme

Following Baines and Palmer (1990), the subgrid-scale orography over one grid-point region is represented by four 
parameters \( \mu \), \( \gamma \), \( \sigma \) and \( \theta \) which stand for the standard deviation, the anisotropy, the slope and the geographical 
orientation of the orography, respectively. These four parameters have been calculated from the US Navy (USN) 
\((10' \times 10')\) data-set.

The scheme uses values of low-level wind velocity and static stability which are partitioned into two parts. The first 
part corresponds to the incident flow which passes over the mountain top, and is evaluated by averaging the wind, 
the Brunt–Väisälä frequency and the fluid density between \( Z_{\text{crit}} \) and \( Z_{\text{H}} \) above the model mean orography. Following 
Wallace et al. (1983), \( 2\mu \) is interpreted as the envelope of the subgrid-scale mountain peaks above the model 
orography. The wind, the Brunt–Väisälä frequency and the density of this part of the low-level flow will be labelled 
\( U_H \), \( N_H \) and \( p_H \), respectively. The second part is the ‘blocked’ flow, and its evaluation is based on a very simple 
interpretation of the non-dimensional mountain height \( H_n \). To first order in the mountain amplitude, the obstacle 
excites a wave, and the sign of the vertical displacement of a fluid parcel is controlled by the wave phase. If a fluid 
parcel ascends the upstream mountain flank over a height large enough to significantly modify the wave phase, its 
vertical displacement can become zero, and it will not cross the mountain summit. In this case the blocking height, 
\( Z_{\text{blk}} \), is the highest level located below the mountain top for which the phase change between \( Z_{\text{blk}} \) and the 
mountain top exceeds a critical value \( H_{n\text{blk}} \), i.e.

\[
\int_{Z_{\text{blk}}}^{H_n} \frac{N}{U_p} \, dz \geq H_{n\text{blk}} \tag{4.9}
\]

In the inequality (4.9), the wind speed, \( U_p(z) \), is calculated by resolving the wind, \( U(z) \), in the direction of the 
flow \( U_H \). Then, if the flow veers or backs with height, (4.9) will be satisfied when the flow becomes normal to
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$U_H$. Levels below this ‘critical’ altitude define the low-level blocked flow. The inequality (4.9) will also be satisfied below inversion layers, where the parameter $N$ is very large. These two properties allow the new parametrization scheme to mimic the vortex shedding observed when pronounced inversions occur (Eling 1989). The upper limit in the equality (4.9) was chosen to be $3\mu$, which is above the subgrid-scale mountain tops. This ensures that the integration in equality (4.9) does not lead to an underestimation of $Z_{blk}$, which can occur because of the limited vertical resolution when using $2\mu$ as an upper limit (a better representation of the peak height), but this upper limit could be relaxed given better vertical resolution.

In the following subsection the drag amplitudes will be estimated combining formulae valid for elliptical mountains with real orographic data. Considerable simplifications are implied and the calculations are, virtually, scale analyses relating the various amplitudes to the sub-grid parameters.

4.2.1 Blocked-flow drag

Within a given layer located below the blocking level $Z_{blk}$, the drag is given by Eq. (4.5). At a given altitude $z$, the intersection between the mountain and the layer approximates to an ellipse of eccentricity

$$ (a', b') = (a, b) \left( \frac{Z_{blk} - z}{z + \mu} \right)^{1/2}, \quad (4.10) $$

where, by comparison with Eq. (4.6), it is also supposed that the level $z = 0$ (i.e. the model mean orography) is at an altitude $\mu$ above the mountain valleys. If the flow direction is taken into account, the length $l(z)$ can be written approximately as

$$ l(z) = 2 \max \left( b \cos \psi, a \sin \psi \right) \left( \frac{Z_{blk} - z}{z + \mu} \right)^{1/2} \quad (4.11) $$

where $\psi$ is the angle between the incident flow direction and the normal ridge direction, $\theta$. For one grid-point region and for uniformly distributed subgrid-scale orography, the incident flow encounters $L/(2a)$ obstacles is normal to the ridge ($\psi = 0$), whereas if it is parallel to the ridge ($\psi = \pi/2$) it encounters $L/(2b)$ obstacles, where $L$ is the length scale of the grid-point region. If we sum up these contributions, the dependence of Eq. (4.11) on $a$ and $b$ can be neglected, and the length $l(z)$ becomes

$$ l(z) = L \left( \frac{Z_{blk} - z}{z + \mu} \right)^{1/2}. \quad (4.12) $$

Furthermore, the number of consecutive ridges (i.e. located one after the other in the direction of the flow) depends on the obstacle shape: there are approximately $L/(2b)$ successive obstacles when the flow is along the ridge, and $L/(2a)$ when it is normal to the ridge. If we take this into account, together with the flow direction, then

$$ l(z) = \frac{L^2}{2} \left( \frac{Z_{blk} - z}{z + \mu} \right)^{1/2} \max \left( \frac{\cos \psi}{a}, \frac{\sin \psi}{b} \right). \quad (4.13) $$

Relating the parameters $a$ and $b$ to the subgrid-scale orography parameters $a = \mu/\sigma$ and $a/b = \gamma$ and, allowing the drag coefficient to vary with the aspect ratio of the obstacle as seen by the incident flow, we have
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\[ r = \frac{\cos^2 \psi + \gamma \sin^2 \psi}{\gamma \cos^2 \psi + \sin^2 \psi}, \]  
(4.14)

and the drag per unit area and per unit height can be written

\[ D_{blk}(z) = -C_d \max(2, 0) \left[ \frac{\sigma}{2 \mu} \left( \frac{Z_{blk} - z}{z + \mu} \right)^{1/2} \max(\cos \psi, \gamma \sin \psi) \frac{U|U|}{2} \right]. \]  
(4.15)

The drag coefficient is modulated by the aspect ratio of the obstacle to account for the fact that \( C_d \) is twice as large for flow normal to an elongated obstacle as it is for flow round an isotropic obstacle. The drag tends to zero when the flow is nearly along a long ridge because flow separation is not expected to occur for a configuration of that kind. It can be shown that the term \( \max(\cos \psi, \gamma \sin \psi) \) is similar to a later form used for the directional dependence of the gravity-wave stress. For simplicity, this later form has been adopted, i.e.

\[ D_{blk}(z) = C_d \max(2, 0) \left[ \frac{\sigma}{2 \mu} \left( \frac{Z_{blk} - z}{z + \mu} \right)^{1/2} \left( B \cos^2 \psi + C \sin^2 \psi \right) \frac{U|U|}{2} \right] \]  
(4.16)

where the constants \( B(\gamma) \) and \( C(\gamma) \) are defined below. The difference between Eq. (4.15) and Eq. (4.16) has been shown to have only a negligible impact on all aspects of the model’s behaviour.

In practice, Eq. (4.16) is suitably resolved and applied to the component from of the horizontal momentum equations. This equation is applied level by level below \( \mu \) and, to ensure numerical stability, a quasi-implicit treatment is adopted whereby the wind velocity \( U \) in Eq. (4.16) is evaluated at the updated time \( t + dt \), while the wind amplitude, \( |U| \), is evaluated at the previous time step.

4.2.2 Gravity-wave drag

This gravity-wave part of the scheme is based on the work of Miller et al. (1989) and Baines and Palmer (1990), and takes into account some three-dimensional effects in the wave stress amplitude and orientation. For clarity and convenience, a brief description is given here. On the assumption that the subgrid-scale orography has the shape of one single elliptical mountain, the mountain wave stress can be written as (Phillips 1984)

\[ (\tau_1, \tau_2) = \rho_H U_H N_H H^2 b G (B \cos^2 \psi_H + C \sin^2 \psi_H, (B - C) \sin \psi \cos \psi_H) \]  
(4.17)

where \( B = 1 - 0.18 \gamma - 0.04 \gamma^2 \), \( C = 0.48 \gamma + 0.3 \gamma^2 \) and \( G \) is a constant of order unity. Furthermore, when \( b \) or \( a \) are significantly smaller than the length \( L \), characteristic of the gridpoint region size, there are, typically, \( L^2/(4ab) \) ridges inside the grid-point region. Summing all the associated forces we find the stress per unit area, viz.

\[ (\tau_1, \tau_2) = \rho_H U_H N_H H^2 \sigma G (B \cos^2 \psi_H + C \sin^2 \psi_H, (B - C) \sin \psi \cos \psi_H) \]  
(4.18)

where \( H \) has been replaced by \( 2 \mu \), and \( a \) by \( \mu/\sigma \).

It is worth noting that, since the basic parameters \( \rho_H, U_H, N_H \) are evaluated for the layer between \( \mu \) and \( 2 \mu \) above the mean orography that defines the model’s lower boundary, there will be much less diurnal cycle in the stress than in previous formulations that used the lowest model levels for this evaluation. The vertical distribution of the gravity-wave stress will determine the levels at which the waves break and slow down the synoptic flow.
Since this part of the scheme is active only above the blocked flow, this stress is now constant from the bottom model level to the top of the blocked flow, $Z_{blk}$. Above $Z_{blk}$, up to the top of the model, the stress is constant until the waves break. This occurs when the total Richardson number, $Ri$, falls below a critical value $Ri_{crit}$, which is of order unity. When the non-dimensional mountain height is close to unity, this algorithm will usually predict wave breaking at relatively low levels; this is not surprising since the linear theory of mountain gravity waves predicts low-level breaking waves at large non-dimensional mountain heights (Miles and Huppert 1969). In reality, the depth over which gravity-wave breaking occurs is more likely to be related to the vertical wavelength of the waves. For this reason, when low-level wave breaking occurs in the scheme, the corresponding drag is distributed (above the blocked flow), over a layer of thickness $\Delta z$, equal to a quarter of the vertical wavelengths of the waves, i.e.

$$
\int_{Z_{fin}}^{Z_{ini}} \frac{N^2}{U_p} \, dz = \frac{\pi}{2}
$$

Above the height $Z_{blk} + \Delta z$ are waves with an amplitude such that $Ri > Ri_{crit}$.

### 4.3 Specification of Subgrid-Scale Orography

For completeness, the following describes how the subgrid-scale orography fields were computed by Baines and Palmer (1990). The mean topographic height above mean sea level over the gridpoint region (GPR) is denoted by $\bar{h}$, and the coordinate $z$ denotes elevation above this level. Then the topography relative to this height $h(x, y) - \bar{h}$ is represented by four parameters, as follows

(i) The net variance, or standard deviation, $\mu$, of $h(x, y)$ in the grid-point region. This is calculated from the US Navy data-set, or equivalent, as described by Wallace et al. (1983). The quantity $\mu$ gives a measure of the amplitude and $2\mu$ approximates the physical envelope of the peaks.

(ii) A parameter $\gamma$ which characterizes the anisotropy of the topography within the grid-point region.

(iii) An angle $\psi$, which denotes the angle between the direction of the low-level wind and that of the principal axis of the topography.

(iv) A parameter $\sigma$ which represents the mean slope within the grid-point region.

The parameters $\gamma$ and $\psi$ may be defined from the topographic gradient correlation tensor

$$
H_{ij} = \frac{\partial h}{\partial x_i} \frac{\partial h}{\partial x_j},
$$

where $x_1 = x$, and $x_2 = y$, and where the terms be calculated (from the USN data-set) by using all relevant pairs of adjacent gridpoints within the grid-point region. This symmetric tensor may be diagonalized to find the directions of the principal axes and the degree of anisotropy. If

$$
K = \frac{1}{2} \left\{ \left( \frac{\partial h}{\partial x} \right)^2 + \left( \frac{\partial h}{\partial y} \right)^2 \right\}, \quad L = \left\{ \left( \frac{\partial h}{\partial x} \right)^2 + \left( \frac{\partial h}{\partial y} \right)^2 \right\} \quad \text{and} \quad M = \frac{\partial h}{\partial x} \frac{\partial h}{\partial y},
$$

the principal axis of $H_{ij}$ is oriented at an angle $\theta$ to the $x$-axis, where $\theta$ is given by

$$
\theta = \frac{1}{2} \arctan(M/L).
$$
This gives the direction where the topographic variations, as measured by the mean-square gradient, are largest. The corresponding direction for minimum variation is at right angles to this. Changing coordinates to \( x', y' \) which are oriented along the principal axes \( x' = x \cos \theta + y \sin \theta \) and \( y' = y \cos \theta - x \sin \theta \), the new values of \( K, L \) and \( M \) relative to these axes, denoted \( K', L' \) and \( M' \), are given by

\[
K' = K, \quad L' = (L^2 + M^2)^{1/2}, \quad M' = 0,
\]

where \( K, L \) and \( M \) are given by Eq. (4.20). The anisotropy of the orography or ‘aspect ratio’. \( \gamma \) is then defined by the equations

\[
\gamma^2 = \left( \frac{\partial h}{\partial y'} \right)^2 \left( \frac{\partial h}{\partial x'} \right)^2
\]

\[
\gamma^2 = \frac{K' - L'}{K' + L'} = \frac{K - (L^2 + M^2)^{1/2}}{K + (L^2 + M^2)^{1/2}}
\]

(4.22)

If the low-level wind vector is directed at an angle \( \psi \) to the \( x \)-axis, then the angle \( \psi \) is given by

\[
\psi = \theta - \psi.
\]

(4.23)

The slope parameter, \( \sigma \), is defined as

\[
\sigma^2 = \left( \frac{\partial h}{\partial x} \right)^2,
\]

(4.24)

i.e. the mean-square gradient along the principal axis.

4.4 CODE

The code mirrors the basic form of the scheme. Hence there is a routine defining all the basic input values for the evaluations of drag, wave stress etc.; a routine to calculate the vertical distribution of wave stress; and a principal routine which computes the wave stress at the surface and the total momentum tendencies, including that from the low-level drag.

The orography parametrization is called from CALLPAR as GWDRAG which in turn calls GWSETUP, and GWPROFIL.

4.4.1 GWSETUP

This defines various reference model levels for controlling the vertical structure of the calculations, and sets up a number of derived atmospheric variables and geometric calculations required to run the scheme:

(a) The definition of the Brunt–Väisälä frequency on half levels

\[
N_{k-1/2}^2 = \frac{2\theta^2}{c_{p \text{dry}}(T^*; T_{k} + T_{k-1})} \left[ 1 - c_{p \text{dry}} \frac{(T_{k} - T_{k-1})}{(p_{k} - p_{k-1})} \right]
\]

(4.25)
(b) The definition of the mean wind components in the layer \( \mu < z < 2\mu \) (where \( \mu \) is the standard deviation of the subgridscale orographic height)

\[
U_{\text{LOW}} = \frac{\sum_{k=\mu}^{k=2\mu} U_k \Delta p_k}{\sum \Delta p_k} \tag{4.26}
\]

and similarly for \( V_{\text{LOW}} \); likewise the mean static stability, \( N_{\text{LOW}} \), and the mean density, \( \rho_{\text{LOW}} \) are calculated.

(c) The calculation of necessary geometry pertaining to geographical orientation of subgridscale orography and wind direction,

\[
\varphi_k = \tan^{-1} \left( \frac{V_k}{U_k} \right) \tag{4.27}
\]

\[
\psi_k = \theta - \varphi_k \tag{4.28}
\]

\[
\bar{\varphi} = \tan^{-1} \left( \frac{V_{\text{LOW}}}{U_{\text{LOW}}} \right) \tag{4.29}
\]

and \( \bar{\psi} = \theta - \bar{\varphi} \), where \( \theta \) is the orientation of ridges relative to east, and the calculation of Phillips (1984) parameters

\[
B = 1 - 0.18\gamma - 0.04\gamma^2, \quad C = 0.48\gamma + 0.3\gamma^2, \tag{4.30}
\]

where \( \gamma \) is the anisotropy of the subgridscale orography.

(d) The calculation of the vertical wind-profile in the plane of the gravity wave stress. Defining

\[
\hat{U}_k = \frac{U_{\text{LOW}}}{|V_{\text{LOW}}|} U_k + \frac{V_{\text{LOW}}}{|V_{\text{LOW}}|} V_k
\]

and similarly for \( \hat{V}_k \), where \( V_{\text{LOW}} = (U_{\text{LOW}} \cdot V_{\text{LOW}}) \), then the wind profile is defined level-by-level as

\[
V^G_k = (\hat{U}_k D_1 + \hat{V}_k D_2) \left( D_1^2 + D_2^2 \right)^{1/2}, \tag{4.31}
\]

where \( D_1 = B - (B - C) \sin^2 \bar{\psi} \) and \( D_2 = (B - C) \sin \bar{\psi} \cos \bar{\psi} \); the values of \( V^G_k \) are also used to compute half level values \( V^G_{k+1/2} \), etc. by linear interpolation in pressure.

(e) The calculation of basic flow Richardson Number

\[
\overline{Ri}_{k-1/2} = N_{k-1/2}^2 \left\{ \frac{p_k - p_{k-1}}{g p_{k-1/2} (V^G_k - V^G_{k-1})} \right\}^2
\]
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(f) The calculation of the depth of the layer treated as ‘blocked’ (i.e. experiencing a direct drag-force due to the subgrid-scale orography). This is given by the value of \( Z_{\text{blk}} \) that is the solution to the finite-difference form of the equation

\[
\int_{Z_{\text{ref}}}^{Z_{\text{blk}}} \frac{N_k}{U_k} \, dz \geq F_c \tag{4.32}
\]

where \( F_c \) is a constant defined later.

(g) The calculation of the layer in which low-level wave-breaking occurs (i.e. the layer experiencing gravity wave breaking (if any) immediately above the ‘blocked’ layer). This is given by the value of \( \Delta z \) that is the solution to the finite difference form of the equation

\[
\int_{Z_{\text{blk}}}^{Z_{\text{blk}} + \Delta z} \frac{N_k}{U_k} \, dz = \frac{\pi}{2} ;
\tag{4.33}
\]

the value of \((Z_{\text{blk}} + \Delta z)\) is not allowed to be less than \(4\mu\).

(h) The calculation of the assumed vertical profile of the subgrid-scale orography needed for the ‘blocking’ computations

\[
z_{\text{DEP}}^k = \frac{Z_{\text{blk}} - z_k}{\eta \left( \frac{z_k}{\mu} \right)} \tag{4.34}
\]

4.4.2 GWPROFIL

This computes the vertical profile of gravity-wave stress by constructing a local wave Richardson number which attempts to describe the onset of turbulence due to the gravity waves becoming convectively unstable or encountering critical layers. This wave Richardson number can be written in the form

\[
\tilde{R}_i = R_i \left( 1 - \frac{1 - \alpha}{\left( 1 + R_i^{1/2} \alpha \right)^2} \right)^{1/2},
\]

where \( R_i \) is the Richardson number of the basic flow. The parameter \( \alpha = N|\delta z|/U_z \) in which \( |\delta z| \) represents the amplitude of the wave and \( U_z \) is the wind speed resolved in the direction of \( \tau_{\text{wave}} \). By requiring that \( \tilde{R}_i \) never falls below a critical value \( \tilde{R}_{i\text{crit}} \) (currently equal to 0.25), values of wave stress are defined progressively from the top of the blocked layer upwards.

When low-level breaking occurs the relevant depth is assumed to be related to the vertical wavelength. Hence a linear (in pressure) decrease of stress is included over a depth \( \Delta z \) given by the solution of Eq. (4.32). The linear decrease of stress is written as

\[
\tau(p_k) = \tau_{Z_{\text{blk}}} + \left( \tau_{Z_{\text{blk}}} - \tau_{Z_{\text{blk}}} \right) \frac{(p_k - p_{Z_{\text{blk}}})}{(p_* - p_{Z_{\text{blk}}})} \tag{4.35}
\]

where the asterisk subscript indicates that the value is at the level \((Z_{\text{blk}} + \Delta z)\).
4.4.3 GWDRAG

This is the main routine. After calling GWSETUP, it defines the gravity-wave stress amplitude in the form,

$$\tau_{\text{wave}} = k p_{\text{LOW}} \sigma \mu (U_{\text{LOW}}^2 + V_{\text{LOW}}^2)^{1/2} (D_1^2 + D_2^2)^{1/2} N_{\text{LOW}}$$

(4.36)

(where $k$ is a constant defined later and $\sigma$ is the mean slope of the subgrid-scale orography) and then calls GWPROCIL. The tendencies due to the wave stresses are then calculated in the form gravity-wave stress amplitude in the form,

$$\tau_{\text{wave}} = k p_{\text{LOW}} \sigma \mu (U_{\text{LOW}}^2 + V_{\text{LOW}}^2)^{1/2} (D_1^2 + D_2^2)^{1/2} N_{\text{LOW}}$$

(4.37)

where $k$ is a constant defined later and $\sigma$ is the mean slope of the subgrid-scale orography.

$$\left( \frac{\partial u}{\partial t} \right)_{\text{wave}} = -g \frac{(\tau_{k+1} - \tau_k)}{(p_{k+1} - p_k)} f(\psi)$$

(4.38)

where $f(\psi)$ is the necessary geometric function to generate components, (similarly for $(\partial v/\partial t)_{\text{wave}}$).

Next the low-level blocking calculations are carried out for levels below $z_h$. These are done level-by-level as follows. Writing the low-level deceleration in the form

$$\left( \frac{\partial u}{\partial t} \right)_{\text{blk}} = -C_d \max \left( 2 - \frac{1}{r}, 0 \right) \frac{\sigma}{2 \mu} \frac{Z_{\text{blk}} - \bar{z}}{\bar{z} + \mu} (B \cos^2 \psi + C \sin^2 \psi) \frac{U[U]}{2},$$

(4.39)

where $r = (\cos^2 \psi + \gamma \sin^2 \psi)/(\gamma \cos^2 \psi + \sin^2 \psi)$ and $B$ and $C$ have been defined earlier, Eq. (4.39) is evaluated in the following partially implicit manner by writing it in the form

$$\left( \frac{\partial u}{\partial t} \right)_{\text{blk}} = \frac{\tilde{U}^{n+1} - U^{n-1}}{2 \Delta t} = -A |U^{n-1}| \tilde{U}^{n+1}$$

then $\tilde{U}^{n+1} = U^{n-1}/(1 + \beta)$ and $\beta = A |U^{n-1}|/2 \Delta t$. Hence

$$\left( \frac{\partial u}{\partial t} \right)_{\text{blk}} = - \left( \beta \right) \frac{U^{n-1}}{2 \Delta t}$$

This calculation is done level-by-level.

Finally the tendencies are incremented. This includes local dissipation heating in the form

$$\left( \frac{\partial T}{\partial t} \right) = \frac{1}{c_p} \frac{\text{DISS}}{2 \Delta t}$$

where $\text{DISS} = \frac{1}{2} \left| \left( U^{n-1} \right)^2 + (V^{n-1})^2 - \left( \tilde{U} \right)^2 \right|$ and $\tilde{U} = U^{n-1} + 2 \Delta t \left( \frac{\partial u}{\partial t} \right)_{\text{sgsor}}$ etc.
APPENDIX A LIST OF SYMBOLS

\( a \)  
- half mountain width in \( x \)-direction

\( B \)  
- function of the mountain anisotropy

\( b \)  
- half mountain width in \( y \)-direction

\( C_d \)  
- drag coefficient

\( D_{blk} \)  
- drag due to flow in blocked layer

\( G \)  
- function of the mountain sharpness

\( H \)  
- maximum mountain height (\( = 2\mu \))

\( h(x,y) \)  
- mountain height profile

\( H_n \)  
- non-dimensional mountain height (\( = NH/|U| \))

\( H_{n_{\text{crit}}} \)  
- critical non-dimensional mountain height

\( L \)  
- length scale of the grid-point region

\( l(z) \)  
- horizontal width of mountain seen by the upstream flow

\( N \)  
- Brunt–Väisälä frequency

\( N_H \)  
- Brunt–Väisälä frequency of un-blocked flow evaluated at height \( H(= 2\mu) \)

\( R_i \)  
- Richardson number

\( R_{i_{\text{crit}}} \)  
- critical Richardson number

\( U \)  
- wind speed in \( x \)-direction

\( U_H \)  
- wind speed of incident un-blocked flow evaluated at height \( H(= 2\mu) \)

\( U_p \)  
- component of the wind speed in the direction of \( U_H \)

\( U_\tau \)  
- component of wind speed in the direction of the stress \( \tau \)

\( V \)  
- wind speed in \( y \)-direction

\( Z_{blk} \)  
- depth of blocked layer

\( \gamma \)  
- anisotropy of the orography (\( = a/b \leq 1 \))

\( \theta \)  
- orientation of the orography

\( \mu \)  
- standard deviation of orography

\( \rho_0 \)  
- density of air at the surface

\( \rho_H \)  
- density of the un-blocked flow evaluated at height \( H(= 2\mu) \)

\( \sigma \)  
- slope of the orography

\( \tau_{blk} \)  
- stress due to blocked flow

\( \tau_{\text{wave}} \)  
- surface stress due to gravity waves

\( \psi \)  
- angle between incident flow and orographic principal axis
Part IV: ‘Physical processes’
CHAPTER 5  Convection

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5.1 INTRODUCTION
Cumulus convection is parametrized by a bulk mass flux scheme which was originally described in Tiedtke (1989). The scheme considers deep, shallow and mid-level convection. Clouds are represented by a single pair of entraining/detraining plumes which describes updraught and downdraught processes. Momentum transport by convective is also included.
5.2 LARGE-SCALE BUDGET EQUATIONS

The contributions from cumulus convection to the large-scale budget equations of heat moisture and momentum are

\[
\left\{ \begin{align*}
\frac{\partial s}{\partial t}_{\text{cu}} &= -\frac{1}{\rho} \frac{\partial}{\partial z} \left[ M_{\text{up}} s_{\text{up}} + M_{\text{down}} s_{\text{down}} - (M_{\text{up}} + M_{\text{down}}) \bar{s} \right] \\
& \quad + L (e_{\text{up}} - e_{\text{down}} - \bar{e}_{\text{subcl}}) - (L_{\text{subl}} - L_{\text{vap}}) (M - \bar{M}) \\
\frac{\partial q}{\partial t}_{\text{cu}} &= -\frac{1}{\rho} \frac{\partial}{\partial z} \left[ M_{\text{up}} q_{\text{up}} + M_{\text{down}} q_{\text{down}} - (M_{\text{up}} + M_{\text{down}}) \bar{q} \right] \\
& \quad - (c_{\text{up}} - e_{\text{down}} - \bar{e}_{\text{subcl}}) \\
\frac{\partial \bar{u}}{\partial t}_{\text{cu}} &= -\frac{1}{\rho} \frac{\partial}{\partial z} \left[ M_{\text{up}} u_{\text{up}} + M_{\text{down}} u_{\text{down}} - (M_{\text{up}} + M_{\text{down}}) \bar{u} \right] \\
\frac{\partial \bar{v}}{\partial t}_{\text{cu}} &= -\frac{1}{\rho} \frac{\partial}{\partial z} \left[ M_{\text{up}} v_{\text{up}} + M_{\text{down}} v_{\text{down}} - (M_{\text{up}} + M_{\text{down}}) \bar{v} \right]
\end{align*} \right. \tag{5.1}
\]

where \( M_{\text{up}}, M_{\text{down}} \) are the net contributions from all clouds to the updraught and downdraught mass fluxes, \( c_{\text{up}} \) and \( e_{\text{down}} \) are the condensation/sublimation in the updraughts, and the evaporation in the downdraughts. \( s_{\text{up}}, s_{\text{down}}, q_{\text{up}}, q_{\text{down}}, u_{\text{up}}, u_{\text{down}}, v_{\text{up}} \) and \( v_{\text{down}} \) are the weighted averages of the dry static energy (\( s \)), the specific humidity (\( q \)), and the horizontal wind components (\( u \) and \( v \)) from all updraughts and downdraughts within a grid box (although individual convective elements are not considered) obtained from the bulk cloud model described below. \( L_{\text{subl}} \) and \( L_{\text{vap}} \) are latent heats of sublimation and vaporization, and \( L \) is the effective latent heat for an ice-water mix (an empirical function of temperature). \( \bar{e}_{\text{subcl}} \) is the evaporation of precipitation in the unsaturated sub-cloud layer, \( \bar{M} \) is the melting of snow and \( \bar{\varphi} \) is the freezing of condensate in the convective updraught. In addition to (5.1) we consider the equations for precipitation

\[
P^{\text{rain}}(z) = \int (G^{\text{rain}} - e_{\text{rain}} - \bar{e}_{\text{subcl}} + \bar{M}) \bar{p} dz , \quad P^{\text{snow}}(z) = \int (G^{\text{snow}} - e_{\text{snow}} - \bar{e}_{\text{subcl}} - \bar{M}) \bar{p} dz \tag{5.2}
\]

where \( P^{\text{rain}}(z) \) and \( P^{\text{snow}}(z) \) are the fluxes of precipitation in the forms of rain and snow at height \( z \). \( G^{\text{rain}} \) and \( G^{\text{snow}} \) are the conversion rates from cloud water into rain and cloud ice into snow. The evaporation of precipitation in the downdraughts \( e_{\text{down}} \), and below cloud base \( \bar{e}_{\text{subcl}} \), have been split into water and ice components, \( e_{\text{rain}}^{\text{water}}, e_{\text{rain}}^{\text{ice}}, e_{\text{snow}}^{\text{water}}, e_{\text{snow}}^{\text{ice}}, \bar{e}_{\text{subcl}}^{\text{water}}, \bar{e}_{\text{subcl}}^{\text{ice}} \).

5.3 CLOUD MODEL EQUATIONS

5.3.1 Updraughts

The updraught of the cloud ensemble is assumed to be in a steady state. Then the bulk equations for mass, heat, moisture, cloud water content and momentum are
Simpson requires knowledge of the cloud-base mass flux and of the mass entrainment and detrainment rates. Cloud-base mass flux is determined for the various types of convection from the closure assumptions discussed in Section 5.4.

Entrainment of mass into convective plumes is assumed to occur (1) through turbulence exchange of mass through the cloud edges and (2) through organized inflow and detrainment is assumed to occur (1) through turbulent exchange and (2) through organized outflow at cloud top. The superscripts (1) and (2) are used to denote the components of the entrainment and detrainment due to turbulent and organized exchanges, respectively:

$$E_{up} = E^{(1)}_{up} + E^{(2)}_{up}, \quad D_{up} = D^{(1)}_{up} + D^{(2)}_{up}$$  \hspace{1cm} (5.4)

5.3.1 (a) Entrainment and detrainment rates. Turbulent entrainment and detrainment are parametrized as

$$E^{(1)}_{up} = \varepsilon^{(1)}_{up} M_{up} \quad D^{(1)}_{up} = \delta^{(1)}_{up} M_{up}$$  \hspace{1cm} (5.5)

where the fractional entrainment/detrainment rates depend inversely on cloud radii in the updraughts ($R_{up}$) (Simpson and Wiggert, 1969; Simpson, 1971):

$$\varepsilon^{(1)}_{up} = \frac{0.2}{R_{up}}, \quad \delta^{(1)}_{up} = \frac{0.2}{R_{up}}$$  \hspace{1cm} (5.6)

By assuming typical cloud sizes for the various types of convection, average values of entrainment/detrainment rates are defined; deep convection is assumed to have a larger radius and so a smaller entrainment rates than shallow convection. In order to keep the scheme simple we use fixed values of turbulent entrainment/detrainment rates for each of the various types of convection:

$$\varepsilon^{(1)}_{up} = \delta^{(1)}_{up} = \begin{cases} 1 \times 10^{-4} \text{ m}^{-1} & \text{for penetrative and midlevel convergence in the presence of large-scale flow convergence} \\ 3 \times 10^{-4} \text{ m}^{-1} & \text{for shallow convection in suppressed conditions} \end{cases}$$  \hspace{1cm} (5.7)
For penetrative convection and mid-level convection we deliberately impose a very small value typical for tropical thunder clouds (Simpson, 1971) so as not to inhibit the penetration of clouds to large heights. For shallow convection we use a value typical for the larger trade wind cumuli (Nitta, 1975), noting that small clouds with much larger entrainment/detrainment rates which detrain immediately above cloud base are not represented in our parametrization. In order to take into account enhanced turbulence in the lower part of the clouds, \( \varepsilon_{\text{up}} \) and \( \delta_{\text{up}} \) are increased in the lowest 150hPa of the cloud in the case of deep and shallow convection. The enhancement factor varies linearly from 4 at cloud base to 1 at 150hPa above cloud base. Turbulent entrainment is only applied over the lowest half of the cloud layer.

### 5.3.1 (b) Organized entrainment and detrainment

Organized entrainment is applied to deep and mid-level convection. The formulation used is discussed in Subsection 5.4.1 below.

Organized detrainment is estimated from the vertical variation of the updraught vertical velocity \( u_{\text{up}} \), which is estimated from the budget equation for the updraught kinetic energy

\[
\frac{\partial K_{\text{up}}}{\partial z} = -\frac{\mu_{\text{up}}}{M_{\text{up}}} (1 + \beta C_d) 2K_{\text{up}} + \frac{1}{f(1 + \gamma)} g \frac{T_{v,\text{up}} - \overline{T}_v}{\overline{T}_v} \tag{5.8}
\]

with

\[
K_{\text{up}} = \frac{u_{\text{up}}^2}{2} \tag{5.9}
\]

where \( K_{\text{up}} \) is the updraught kinetic energy, \( T_{v,\text{up}} \) is the virtual temperature of the updraught and \( \overline{T}_v \) the virtual temperature of the environment. \( \mu_{\text{up}} \) is a mixing coefficient which is equal to the entrainment mass flux \( (E_{\text{up}}) \), or the detrainment mass flux \( (D_{\text{up}}) \) if this is larger. As entrainment is set to zero in the upper part of the cloud layer, use of detrainment mass flux in this region better represents the effect of mixing and vertical pressure gradients in the upper part of deep convective clouds, reducing vertical velocity and reducing overshoot of convective towers into the lower stratosphere.

\( \gamma \) (= 0.5) is the virtual mass coefficient (Simpson and Wiggert 1969), the factor \( f \) (= 2) is introduced because the flow is highly turbulent (Cheng et al. 1980) and for the drag coefficient a value of \( C_d = 0.506 \) is used (Simpson and Wiggert 1969). The value for \( \beta \) is 1.875. The cloud base value of the updraught velocity is chosen as 1 m s\(^{-1}\).

\( u_{\text{up}} \) enters the scheme in several ways: (i) for the generation and fallout of rain (Section 5.6), (ii) to determine penetration above the zero-buoyancy level and the top of cumulus updraughts and (iii) to specify detrainment below the top of the updraught.

\( u_{\text{up}} \) determines the level to which convection penetrates (where it reduced to zero). This allows convection to penetrate above its level of neutral buoyancy. Organized detrainment is estimated by equating the decrease in updraught vertical velocity due to negative buoyancy at the top of the cloud to the decrease in mass flux with height:

\[
\frac{M_{\text{up}}(z)}{M_{\text{up}}(z + \Delta z)} = \sqrt{\frac{K_{\text{up}}(z)}{K_{\text{up}}(z + \Delta z)}} \tag{5.10}
\]

This assumes that the cloud area remains constant in the detraining layer and neglects the vertical variation of buoyancy. Eq. (5.10) defines the reduction of mass flux with height, which combined with the updraught continuity equation (Eq. (5.3)) gives the organised detrainment rate.
5.3.2 Downdraughts

Downdraughts are considered to be associated with convective precipitation from the updraughts and originate from cloud air influenced by the injection of environmental air. Following Fritsch and Chappell (1980) and Foster (1958), the Level of Free Sinking (LFS) is assumed to be the highest model level (below the level of minimum moist static energy) where a mixture of equal parts of cloud and saturated environmental air at the wet-bulb temperature becomes negative buoyant with respect to the environmental air. The downdraught mass flux is assumed to be directly proportional to the upward mass flux. Following Johnson (1976, 1980) the mass flux at the LFS is specified from the updraught mass flux at cloud base as

\[ (M_{\text{down}})_{\text{LFS}} = \eta(M_{\text{up}})_{\text{base}} \quad \text{with} \quad \eta = -0.3 \quad (5.11) \]

The vertical distribution of the downdraught mass flux, dry static energy, moisture and horizontal momentum below the LFS are determined by entraining/detraining plume equations similar to those for the updraught;

\[
\begin{align*}
- \frac{\partial M_{\text{down}}}{\partial z} &= E_{\text{down}} - D_{\text{down}} \\
- \frac{\partial (M_{\text{down}}s_{\text{down}})}{\partial z} &= E_{\text{down}}s - D_{\text{down}}s_{\text{down}} + L\rho e_{\text{down}} \\
- \frac{\partial (M_{\text{down}}q_{\text{down}})}{\partial z} &= E_{\text{down}}q - D_{\text{down}}q_{\text{down}} + \rho e_{\text{down}} \\
- \frac{\partial (M_{\text{down}}u_{\text{down}})}{\partial z} &= E_{\text{down}}u - D_{\text{down}}u_{\text{down}} \\
- \frac{\partial (M_{\text{down}}v_{\text{down}})}{\partial z} &= E_{\text{down}}v - D_{\text{down}}v_{\text{down}}
\end{align*}
\]

\( e_{\text{down}} \) is the evaporation of convective rain to maintain a saturated descent; the moistening and cooling of the environmental air injected at LFS is also due to evaporating rain.

Entrainment and detrainment in downdraughts are highly uncertain as relevant data are not available. As for the updraught, both turbulent and organized entrainment/detrainment are considered.

5.3.2 (a) Turbulent entrainment and detrainment. For turbulent mixing

\[ \varepsilon_{\text{down}}^{(1)} = \delta_{\text{down}}^{(1)} = 2 \times 10^{-4} \text{m}^{-1} \quad (5.13) \]

5.3.2 (b) Organized entrainment and detrainment. Organized entrainment for the downdraught is based upon a formulation suggested by Nordeng (1994);

\[
\varepsilon_{\text{down}}^{(2)} = \frac{\left\{ \frac{T_{\text{v,down}} - T_{\text{down}}r_{\text{down}}}{T_{\text{v}}} \right\}}{z} \quad (5.14)
\]

\[ (w_{\text{down}})^{\text{LFS}} = \frac{1}{2} \int \left\{ \frac{T_{\text{v,down}} - T_{\text{down}}r_{\text{down}} - T_{\text{v}}}{T_{\text{v}}} \right\} dz \]

where \( w_{\text{down}}^{\text{LFS}} \) is the vertical velocity in the downdraught at the LFS (set to 1 m s\(^{-1}\)).
The scheme has no explicit rain water equation for the downdraught and so \( r_{\text{down}} \) is estimated by

\[
r_{\text{down}} = \sum_{k=1}^{n_{\text{lev}}} r_{\text{up}}^k M_{\text{up}}^k
\]

Organized detrainment from the downdraught occurs when either the downdraught becomes positively buoyant or approaches the surface. If the downdraught remains negatively buoyant until it reaches the surface then the mass flux is decreased linearly over the bottom three model levels for L31 and L50 versions of the IFS. In L60 versions the downdraught is detrained over the lowest seven model levels, maintaining an outflow depth of about 50hPa as in lower resolution versions of the model. However if a downdraught becomes positively buoyant during its descent, it is detrained over one level, except where this occurs at cloud base, when the downdraught fluxes are decreased linearly (deep convection) or quadratically (mid-level convection) to zero at the surface.

**5.4 Convective types**

In using a bulk mass flux scheme, as opposed to a scheme which considers an ensemble of convective clouds (such as that of Arakawa and Schubert, 1974), some determination of convective cloud type must be made so that appropriate choices can be made for the cloud properties. Firstly it must be determined if the profile can support convection from the surface layer. If on carrying out an undilute ascent from the surface layer a cloud base is found where the parcel buoyancy is greater than \(-0.5\) K, then either deep or mid-level convection is initiated. If no such cloud base is found then higher levels of the model are tested for a buoyant layer, mid-level convection being initiated from the lowest such level.

For convection initiating from the surface, the original version of the convection scheme (Tiedke 1989) used a comparison of moisture convergence and surface evaporation to determine whether convection was deep or shallow. However this is now done on the basis of the depth of the convective cloud. If the cloud depth exceeds 200 hPa then deep convection is assumed, shallow convection if not.

Once the type of convection has been determined its intensity (controlled by the cloud-base mass flux) is determined as outlined below.

**5.4.1 Deep convection**

Following Frisch and Chappell (1980) and Nordeng (1994), the cloud base mass flux for deep convection is estimated from assuming that convection acts to reduce the convective available potential energy (CAPE) towards zero over a specified time scale \( \tau \):

\[
\frac{\partial \text{CAPE}}{\partial t} = -\frac{\text{CAPE}}{\tau} = \int_{z_{\text{base}}}^{z_{\text{top}}} \frac{\bar{g} \frac{\partial T_v}{\partial z}}{T_v} \mathrm{d}z = \int_{z_{\text{base}}}^{z_{\text{top}}} M_{\text{clw}} \frac{\bar{g} \frac{\partial T_v}{\partial z}}{T_v} \mathrm{d}z
\]

where

\[
M_{\text{clw}} = M_{\text{up}} + M_{\text{down}} = \alpha [M_{\text{up}}]_{\text{base}} + \beta [M_{\text{down}}]_{LFS}
\]

where \( \alpha \) and \( \beta \) describe the vertical variation of the updraught and downdraught mass flux due to entrainment and
detrainment and the subscript ‘base’ refers to cloud-base quantities. As the downdraught mass flux at the LFS is linked to the updraught mass flux at cloud base (Eq. (5.11)) then,

\[ M_{\text{cld}} = [M_{\text{up}}]_{\text{base}} (\alpha + \beta \eta) \]  

(5.18)

Using Eq. (5.18) in Eq. (5.16) results in an expression for the cloud base mass flux. CAPE is estimated from the parcel ascent incorporating the effects of water loading,

\[ \text{CAPE} = \int_{z_{\text{base}}}^{z_{\text{top}}} g \left( \frac{T_{\text{up}} - T_v}{T_v} - l_{\text{up}} \right) dz \]  

(5.19)

In practice the vertical variation of the updraught and downdraught mass fluxes (\( \alpha \) and \( \beta \) above) is estimated from an initial ascent using an arbitrary value for the updraught mass flux at cloud base followed by a downdraught calculation. Using these estimates the updraught mass flux at cloud base is recomputed and downdraught mass fluxes rescaled. A second updraught ascent is then computed to revise the updraught properties.

The adjustment time scale \( \tau \) is rather arbitrary but experience suggests that to prevent grid-scale saturation it must be set such that

\[ M_{\text{cld}} = \rho \overline{w} \]  

(5.20)

where \( \overline{w} \) is the grid-scale vertical velocity. The magnitude of the grid-scale vertical velocity increases with increasing resolution, and therefore the relaxation timescale is also made dependent on model resolution. At resolutions smaller than T159 it is set to 2 hours, from T159 to T319 it is set to 1 hour and at resolutions finer than T319 it is set to 20 minutes.

The vertical distribution of the updraught mass flux above cloud base is determined by assuming that there is organized entrainment which is directly proportional to the large-scale moisture convergence as

\[ E_{\text{up}}^{(2)} = -\frac{\beta}{q} (\nabla \cdot \overline{q} + \overline{w} \frac{\partial q}{\partial z}) \]  

(5.21)

Organized entrainment is only considered in the lower part of the cloud layer where large-scale convergence is encountered, that is, below the level of strongest vertical ascent. The idea to link the cloud mass flux directly to the large-scale moisture convergence has first been advocated as a parametrization by Lindzen (1981) who indicated that it may provide vertical profiles of mass flux and convective heating in good agreement with observations. The assumption (5.21) ensures that the vertical distribution of the convective mass flux follows that of the large-scale ascent which is partly supported by diagnostic studies for tropical convection (e.g. Cheng et al., 1980; Johnson, 1980).

### 5.4.2 Shallow convection

Here we consider cumulus convection, which predominantly occurs in undisturbed flow, that is in the absence of large-scale convergent flow. Typical examples are trade-wind cumuli under a subsidence inversion, convection occurring in the ridge region of tropical easterly waves and daytime convection over land. This type of convection seems to be effectively controlled by sub-cloud layer turbulence. In fact, most of the diagnostic studies carried out
for trade-wind cumuli show that the net upward moisture flux at cloud-base level is nearly equal to the turbulent moisture flux at the surface (LeMone and Pennell, 1976). In regions of cold air flowing over relatively warm oceans the then relatively high sensible heat flux has been found to be of significant importance. We therefore derive the mass flux at cloud base on a balance assumption for the sub-cloud layer based on the moist static energy budget:

\[ [M_{up}(h_{up} - \overline{h})]_{base} = -\int_{0}^{h} \left( \nabla \cdot \nabla \overline{h} + \overline{\rho} \frac{\partial}{\partial z} \overline{h} + c_{p} \left( \frac{\partial}{\partial t} \overline{T} \right) \right)_{rad} + \frac{1}{\overline{\rho}} \frac{\partial}{\partial z} \left( \overline{\rho \overline{w} \overline{h}} \right)_{turb} \overline{\rho} \, dz \]  

(5.22)

with

\[ \overline{h} = c_{p} T + L_{q} + g z \]  

(5.23)

The moisture supply to the shallow cumulus is largely through surface evaporation as the contributions from large-scale convergence are either small or even negative, such as in the undisturbed trades where dry air is transported downward to lower levels.

An initial estimate for the updraft base mass flux is obtained using Eq. (5.22). If downdraughts occur (relatively rare for shallow convection due to the low precipitation rates), then a revised estimate is made accounting for the impact of downdraughts upon the sub-cloud layer, the l.h.s. of Eq. (5.22) being replaced by

\[ [M_{up}(h_{up} - \overline{h})]_{base} + [M_{down}(h_{down} - \overline{h})]_{base} = [M_{up}(h_{up} - \overline{h})]_{base} + [\beta \eta M_{up}(h_{down} - \overline{h})]_{base} \]  

(5.24)

Again downdraught properties are obtained using the original estimate of the updraft base mass flux and then rescaled by the revised value. For the updraft a second ascent is calculated using the revised value of the base mass flux.

No organized entrainment is applied to shallow convection. As turbulent entrainment and detrainment rates are equal, the mass flux remains constant with height until reducing at cloud top by organized detrainment.

5.4.3 Mid-level convection

Mid-level convection, that is, convective cells which have their roots not in the boundary layer but originates at levels above the boundary layer, often occur at rain bands at warm fronts and in the warm sector of extratropical cyclones (Browning et al. 1973; Houze et al. 1976; Herzegh and Hobbs 1980). These cells are probably formed by the lifting of low level air until it becomes saturated (Wexler and Atlas 1959) and the primary moisture source for the clouds is from low-level large-scale convergence (Houze et al. 1976). Often a low-level temperature inversion exists that inhibits convection from starting freely from the surface; therefore convection seems to be initiated by lifting low-level air dynamically to the level of free convection. This occurs often in connection with mesoscale circulations which might be related to conditionally symmetric instability (Bennets and Hoskins 1979; Bennets and Sharp 1982) or a wave-CISK mechanism (Emanuel 1982).

Although it is not clear how significant the organization of convection in mesoscale rain bands is for the large-scale flow, a parametrization should ideally account for both convective and mesoscale circulations. Such a parametrization, however, is presently not available and we must therefore rely on simplified schemes. Here we use a parametrization which in a simple way considers the finding of the diagnostic studies mentioned above. We assume that convection is activated when there is a large-scale ascent, the environmental air is sufficiently moist, i.e., of relative humidity in excess of 80%, and convectively unstable layer exists above (i.e. at cloud base the buoyancy is greater than –0.5 K).
The convective mass flux at cloud base is set equal to the vertical mass transport by the large-scale flow at that level:

\[
\bar{\rho}_{\text{base}} \bar{w}_{\text{base}} = (M_{\text{up}})_{\text{base}} + (M_{\text{down}})_{\text{base}} = (M_{\text{up}})_{\text{base}} + \beta \eta (M_{\text{up}})_{\text{base}} = (M_{\text{up}})_{\text{base}} (1 + \beta \eta)
\]  

(5.25)

following the notation of Subsection 5.4.1 above. Again two estimates of the updraught base mass flux are made; first neglecting downdraughts, followed by a revised estimate if downdraughts occur. The closure ensures that the amount of moisture which is vertically advected through cloud base by the large-scale ascent is fully available for generation of convective cells.

In addition to the injection of mass through cloud base, we assume again that cloud air is produced by moisture convergence above cloud base through organised entrainment in the same way as for penetrative convection as given by (5.21).

### 5.5 Sub-cloud layer

The first level at which convective mass, momentum and thermodynamic fluxes are estimated is cloud base. To represent the effects of convective updraughts on the sub-cloud layer a simple scaling of cloud base fluxes is applied in which they decrease to zero at the surface through the sub-cloud layer.

Care must be taken to ensure that fluxes of liquid water are zero below cloud base. Through the cloud base level an interpolation of fluxes liquid water static energy and total water content is used to estimate fluxes of dry static energy and water vapour mixing ratio in the level immediately below cloud base:

\[
(M_s)_{\text{up}}^{\text{base}+1} = (Z^n)(M_s)_{\text{up}}^{\text{base}} - L(M_l)_{\text{up}}^{\text{base}}
\]

\[
(M_q)_{\text{up}}^{\text{base}+1} = (Z^n)(M_q)_{\text{up}}^{\text{base}} + (M_l)_{\text{up}}^{\text{base}}
\]

\[
(M_l)_{\text{up}}^{\text{base}+1} = 0
\]

(5.26)

where \((\phi)^{\text{base}+1}\) refers to the value of \(\phi\) at the level immediately below cloud base. \(Z\) is given by

\[
Z = \frac{p_{\text{surf}} - p_{\text{base}+1}}{p_{\text{surf}} - p_{\text{base}}}
\]

(5.27)

and \(p_{\text{surf}}\) is the surface pressure.

For deep and shallow convection \(n\) is set to 1 (implying a linear decrease in the flux with pressure below cloud base) while for mid-level convection \(n\) is equal to 2 (implying a quadratic reduction in flux below cloud base).

For the remainder of the sub-cloud layer, fluxes at level ‘B+1’ are reduced to zero at the surface using \((Z^n)\) recomputed as

\[
Z = \frac{p_{\text{surf}} - p_k}{p_{\text{surf}} - p_{\text{base}+1}}
\]

(5.28)

where \(p_k\) is the pressure at level \(k\) of the model.

The cloud-mass and momentum fluxes in the sub-cloud layer are treated in a similar manner.
5.6 CLOUD MICROPHYSICS

5.6.1 Freezing in convective updraughts
We assume that condensate in the convective updraughts freezes in the temperature range \(250.16\text{K} < T < 273.16\text{K}\) maintaining a mixed phase within that range according to (6.6) (see Chapter 6 ‘Clouds and large-scale precipitation’).

5.6.2 Generation of precipitation
The conversion from cloud water/ice to rain/snow is treated in a consistent way with that in the large-scale precipitation scheme by using a formulation following Sundqvist (1978)

\[
\bar{\rho} G^{\text{precip}} = M_{\text{up}}\frac{c_0}{0.75 \bar{w}_{\text{up}}} \left[ l_{\text{up}} \left[ 1 - \exp\left\{ -\left( \frac{l_{\text{up}}}{l_{\text{crit}}} \right)^2 \right\} \right] \right]
\]  

(5.29)

where \(c_0 = 1 \times 10^{-3} \text{s}^{-1}\) and \(l_{\text{crit}} = 0.5 \times 10^{-3} \text{g kg}^{-1}\). \(\bar{w}_{\text{up}}\) is the updraught vertical velocity and is limited to a maximum value of 10 m s\(^{-1}\) in Eq. (5.29). The value of the autoconversion coefficient has been increased from values used in previous cycles of the convection scheme by a factor of 1.5.

Sundqvist (1988) takes account of the Bergeron-Findeisen process for temperatures below \(-5^\circ\text{C}\) through a temperature dependent modification of \(c_0\) and \(l_{\text{crit}}\):

\[
c'_0 = c_0c_{\text{BF}}
\]
\[
l'_\text{crit} = l_{\text{crit}}c_{\text{BF}}
\]

(5.30)

where

\[
c_{\text{BF}} = 1 + 0.5\sqrt{\max\left(\frac{T_{\text{BF}} - T_{\text{up}}}{T_{\text{up}} - T_{\text{ice}}}\right)}\text{ for } T < T_{\text{BF}}
\]
\[
c_{\text{BF}} = 1\text{ for } T > T_{\text{BF}}
\]

(5.31)

\(T_{\text{BF}} = 268.16\text{K}\), and \(T_{\text{ice}} = 260\text{K}\).

Eq. (5.29) is integrated analytically in the vertical.

5.6.3 Fallout of precipitation
The fallout of rain water/snow is parametrized as (e.g. Kuo and Raymond, 1980)

\[
\bar{\rho} S^{\text{fallout}} = M_{\text{up}} \frac{V}{\bar{w}_{\text{up}} \Delta z} r_{\text{up}}
\]

(5.32)

where \(\Delta z\) is the model layer depth. The terminal velocity \(V\) is parametrized as (Liu and Orville, 1969)

\[
V = 5.32 r_{\text{up}}^{0.2}
\]

(5.33)

where \(r_{\text{up}}\) is given in units of g kg\(^{-1}\). Since the fall speed of ice particles is smaller than that of water droplets, only half the value of \(V\) calculated with Eq. (5.27) is used for ice. In estimating the fallout of precipitation in the...
mixed phase region of the cloud a weighted mean of the fall speed for ice and water precipitation is used. Eq. (5.32) is integrated analytically in the vertical

5.6.4 Evaporation of rain

The evaporation of convective rain is parametrized following a proposal of Kessler (1969), where the evaporation is assumed to be proportional to the saturation deficit \( q_{\text{sat}} - q \) and to be dependent on the density of rain \( \rho_{\text{rain}} \) (g m\(^{-3}\))

\[
e^{\text{rain}} = \alpha_1 (q_{\text{sat}} - q) \rho_{\text{rain}}^{13/20}
\]

(5.34)

where \( \alpha_1 \) is a constant being zero for \( q > q_{\text{sat}} \).

As the density of rain \( \rho_{\text{rain}} \) is not given by the model it is convenient to express it in terms of the rain intensity \( R \) (g m\(^{-2}\)s\(^{-1}\)) as

\[
R = \rho_{\text{rain}} V_{\text{rain}}
\]

(5.35)

where \( V_{\text{rain}} \) is the mean fall speed of rain drops which again is parametrized following Kessler (1969).

\[
V_{\text{rain}} = \alpha_2 \rho_{\text{rain}}^{1/8} \sqrt{\frac{\rho_{\text{sur}}}{\rho}}
\]

(5.36)

(Note that this is different from the formulation used in the estimation of the fallout of precipitation).

Thus we have

\[
e^{\text{rain}} = \alpha_1 (q_{\text{sat}} - q) \left\{ \left( \frac{\sqrt{\rho_{\text{sur}}}}{\rho} \right)^{13/20} \right\}^{8/9}
\]

(5.37)

Since the convective rain takes place only over a fraction \( C_{\text{conv}} \) of the grid area, the evaporation rate at level \( k \) becomes

\[
e^{\text{rain}} = C_{\text{conv}} \alpha_1 (q_{\text{sat}} - q) \left\{ \frac{\sqrt{\rho_{\text{sur}}}}{\rho} \frac{R}{C_{\text{conv}}} \right\}^{8/9}
\]

(5.38)

where the constants have the following values (Kessler, 1969)

\[
\alpha_1 = 5.44 \times 10^{-4} \quad \alpha_2 = 5.09 \times 10^{-3} \quad \alpha_3 = 0.5777
\]

In view of the uncertainty of the fractional area of precipitating clouds a constant value of \( C_{\text{conv}} = 0.05 \) is assumed.

The evaporation rate is calculated implicitly in the model by means of

\[
2g \frac{\partial (R^{1/2})}{\partial p} = -A
\]

(5.39)

which follows from
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\[ e^{\text{rain}} = AR^{1/2} \quad A = \alpha_1 (q_{sat} - q) \left( \sqrt{\frac{\rho}{p_{\text{surf}}}} \frac{1}{C_{\text{conv}}} \right)^{1/2} \]  

(5.40)

and

\[ e^{\text{rain}} = \frac{1}{\rho} \frac{\partial R}{\partial z} = -\frac{\partial R}{\partial p} \]  

(5.41)

5.6.5 Melting and freezing of precipitation

Melting of snow falling across the freezing level \((T_0)\) is parametrized by a simple relaxation towards \(T_0\):

\[ M = \frac{\epsilon_p (T - T_0)}{\tau} \]  

(5.42)

where \(M\) is the amount of snow (kg m\(^2\) s) melting and \(\tau\) is a relaxation time scale which decreases with increasing temperature,

\[ \tau = \frac{5 \times 3600}{\{ 1 + 0.5(T - T_0) \}} \]  

(5.43)

The parametrization may produce melting over a deeper layer than observed (Mason 1971) but this has been intentionally introduced to account implicitly for the effects of vertical mixing which may develop in response to the production of negative buoyancy.

5.7 Link to cloud scheme

Before the introduction of the prognostic cloud scheme (see Chapter 6 ‘Clouds and large-scale precipitation’) water detrained from convection \((D_{w}\ell_{up})\) was evaporated instantaneously. However with the prognostic cloud scheme water detrained from convection is taken to be a source of cloud mass increasing the cloud fraction and water content of clouds;

\[ \frac{\partial \alpha}{\partial t} = (1 - \alpha) \frac{D_{w}\ell_{up}}{\rho} \]  

(5.44)

\[ \frac{\partial \bar{I}}{\partial t} = \frac{D_{w}\ell_{up}}{\rho} \]

where \(\alpha\) is the cloud fraction and \(\bar{I}\) the grid-box mean cloud water.

5.8 Momentum transports

Equation set (5.3) includes a treatment of the vertical transport of horizontal momentum by convection. Studies have shown that for deep convection momentum transports are over estimated by the plume models unless the effects of cloud scale horizontal pressure gradients are included (Gregory et al. 1997b). For unorganised convection the effects of the pressure gradients are to adjust the in-cloud winds towards those of the large-scale flow. This can be represented by an enhanced turbulent entrainment rate in the cloud momentum equations. To ensure mass con-
continuity the turbulent detrainment rate is also increased by an equivalent amount. As the air entrained as detrained have differing properties this adjusts the in-cloud wind back towards the large-scale value.

Hence for deep and mid-level convection the turbulent entrainment and detrainment rates used in the updraught momentum equation are

\[ \varepsilon_{up}^{k,(1), (u,v)} = \varepsilon_{up}^{(1)} + \lambda \varepsilon_{up}^{(1)} \]
\[ \delta_{up}^{k,(1), (u,v)} = \delta_{up}^{(1)} + \lambda \delta_{up}^{(1)} \] (5.45)

where \( \phi \) is the value of \( \phi \) in level \( k \), while \( \varepsilon_{up}^{(1)} \) and \( \delta_{up}^{(1)} \) are given by equation (5.7).

When \( \varepsilon_{up}^{k,(1)} > 0 \) (below the mid-level of the cloud) \( \lambda = 2 \), while if \( \varepsilon_{up}^{k,(1)} = 0 \) (in the upper part of the cloud) then \( \lambda = 3 \). Gregory (1997) suggests that the above formulation provides an adequate description of the effects of cloud scale pressure gradients in cases of deep convection. For shallow convection and downdraughts it is assumed that the effects of the pressure gradient term can be neglected and no enhancement of the entrainment rates in the momentum equations is applied. This formulation limits the momentum transports to be downgradient. Up-gradient transports by highly organized convective systems (e.g. African squall lines) are not captured by this method.

The definition of the horizontal wind in the updraught and downdraught at cloud base and LFS is not well known. For the updraught the value at cloud base is set to an average of the large-scale winds over the depth of the sub-cloud layer. For the downdraught the initial values at the LFS are set equal to the average values of the winds in the updraught and those of the large-scale flow.

### 5.9 Discretization of the model equations

The flux divergence in the large-scale budget equations (5.1) and in the cloud equations (5.3) and (5.12) are approximated by centred finite differences as

\[ \frac{1}{p} \frac{\partial (Ma)}{\partial z} = g \left( \frac{M_{k+1/2} \alpha_{k+1/2} - M_{k-1/2} \alpha_{k-1/2}}{p_{k+1/2} - p_{k-1/2}} \right) \] (5.46)

The definition of the large-scale variables at half levels pose a problem, when the half-level values defined by linear interpolation of full-level values very noisy profiles evolve in time particularly with regard to humidity. Much smoother profiles are obtained when the half-level values are determined by downward extrapolation from the next full level above along a cloud-ascent through that level:

\[ \bar{T}_{k+1/2} = \bar{T}_k + \left( \frac{\partial \bar{T}}{\partial p} \right)_{h_{sat}} (p_{k+1/2} - p_k) \]
\[ \bar{q}_{k+1/2} = \bar{q}_k + \left( \frac{\partial \bar{q}}{\partial p} \right)_{h_{sat}} (p_{k+1/2} - p_k) \] (5.47)

where \( h_{sat} = c_p T + g z + L_{sat} \) is the saturation moist static energy. Using an extrapolation like (5.47) for calculating the downward transports is also more consistent with the calculation of the updraughts where cloud air is transported upwards through level \( k + 1/2 \) with the thermal state below that level and equally with the downdraughts which depend only on values of \( s \) and \( q \) above that level. Similarly, because of (5.46) the downward transport of
environmental air through the same level accounts now only for thermal properties above that level. The choice of a moist adiabat for extrapolation is dictated by the property of the moist static energy which is, by convection in the absence of downdraughts, only changed through the fluxes of moist static energy

\[
\left( \frac{\partial h}{\partial t} \right)_{cu} = \frac{1}{\rho} \frac{\partial}{\partial z} [M_{up}(h_{up} - \bar{h})] \tag{5.48}
\]

As the lines of the saturation moist static energy \( h_{sat} \) through point \((p_{h+1/2}, \bar{T}_{h-1/2})\) and the updraught moist static energy are almost parallel, apart from entrainment effects, the difference \( h_{up} - \bar{h} \) is little affected by the vertical discretization.

For horizontal winds, values at model half levels are set to those on the full model level below.

The ascent in the updraughts is obtained by vertical integration of (5.3). Starting at the surface the condensation level (equal to the lowest half-level which is saturated or supersaturated and where buoyancy is greater than \(-0.5 \text{ K}\) ) is determined from an adiabatic ascent. The cloud profile above cloud base is determined layer by layer by first doing a dry adiabatic ascent with entrainment and detrainment included and then adjusting temperature and moisture towards a saturated state, taking into account condensation and freezing processes. The buoyancy of the parcel is calculated taking into account the effects of cloud and precipitation water loading i.e.

\[
B = T_{up}(1 + 0.608q_{up} - l_{up} - r_{up}) - T_s(1 + 0.608q_s) \tag{5.49}
\]

Special care has to be taken in the discretization of (5.8) because of overshooting effects. A centred differencing scheme is used so that

\[
\frac{K_{up,k+1/2} - K_{up,k+1/2}}{z_{k+1/2} - z_{k+1/2}} = \frac{E_{up,k}}{M_{up,k+1/2}} \left( 1 + \beta C_d \right) \{K_{up,k-1/2} + K_{up,k+1/2} \} + \frac{1}{f(1 + \gamma)} \frac{1}{2} \left[ \frac{T_{v,up} - \bar{T}_v}{k-1/2} + \frac{T_{v,up} - \bar{T}_v}{k+1/2} \right] \tag{5.50}
\]

Finally, we mention that for numerical reasons the environmental air must not be convectively unstably stratified:

\[
\bar{s}_{k-1/2} \geq \bar{s}_{k+1/2} \tag{5.51}
\]

In fact, one of the forecasts with the ECMWF global model became numerically unstable when (5.22) was not imposed.

### 5.10 Structure of code

The parametrization of cumulus convection is performed in subroutines shown in Fig. 5.1.

**Figure 5.1** Structure of convection scheme

**CUCALLN**

Provides interface of routines for cumulus parametrization. It takes the input values through arguments from CALLPAR and returns updated tendencies of \( T, q, a, l, u \) and \( v \), as well as convective precipitation rates.
Chapter 5 ‘Convection’

**CUMASTRN**
Master routine for convection scheme.

**CUININ**
Initializes variables for convection scheme (including vertical interpolation to the half model levels).

**CUBASEN**
Calculates condensation level for surface air and sets updraft base variables.

**CUASCN**
Calculates ascent in updraughts. **CUASCN** is called twice, the second time after downdraughts have been calculated taking account of the CAPE adjustment closure for deep convection and downdraughts for mid-level and shallow convection Routines **CUENTR** and **CUBASMCN** are called from **CUASCN**.

**CUENTR**
Calculated entrainment and detrainment rates.

**CUBASMCN**
Calculates cloud base properties of mid-level convection.

**CUDLFSN**
Calculates the level of free sinking for downdraughts.

**CUDDRAFN**
Calculates the downdraught descent.

**CUFLXN**
Calculates final convective fluxes and surface precipitation rates taking into account of melting/freezing and the evaporation of falling precipitation.

**CUDDQN**
Calculates the tendencies of $T$ and $q$ from convection.

**CUUDUV**
Calculates the tendencies of $u$ and $v$ from convection

**CUADTQ**
Calculates super/sub saturation and adjusts $T$ and $q$ accordingly.

**EXTERNALS**
Subroutine SATUR for calculating saturation mixing ratio.

**PARAMETERS**
Defined in subroutine SUCUM called from INIPHY.

### APPENDIX A LIST OF SYMBOLS

<table>
<thead>
<tr>
<th>Symbol</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>CAPE</td>
<td>Convective available potential energy</td>
</tr>
<tr>
<td>$C_d$</td>
<td>Drag coefficient</td>
</tr>
<tr>
<td>$c_{conv}$</td>
<td>Fraction of grid square occupied by convection</td>
</tr>
<tr>
<td>$c_p$</td>
<td>Specific at constant pressure for moist air</td>
</tr>
<tr>
<td>$c_{up}$</td>
<td>Condensation/sublimation in the updraughts</td>
</tr>
<tr>
<td>$D_{up}$</td>
<td>Rate of mass detrainment in the updraughts</td>
</tr>
<tr>
<td>$D_{down}$</td>
<td>Rate of mass detrainment in the downdraughts</td>
</tr>
<tr>
<td>$E_{up}$</td>
<td>Rate of mass entrainment in the updraughts</td>
</tr>
<tr>
<td>$E_{down}$</td>
<td>Rate of mass entrainment in the downdraughts</td>
</tr>
<tr>
<td>$e_{rain}$</td>
<td>Evaporation of rain</td>
</tr>
<tr>
<td>$e_{down}$</td>
<td>Evaporation of precipitation (rain and snow) in the downdraughts</td>
</tr>
<tr>
<td>$e_{rain}$</td>
<td>Evaporation of rain in the downdraughts</td>
</tr>
<tr>
<td>$e_{snow}$</td>
<td>Evaporation of snow in the downdraughts</td>
</tr>
<tr>
<td>$e_{subcld}$</td>
<td>Evaporation of precipitation (rain and snow) in the unsaturated sub-cloud layer</td>
</tr>
<tr>
<td>$e_{subcld}$</td>
<td>Evaporation of rain in the unsaturated sub-cloud layer</td>
</tr>
<tr>
<td>$e_{subcld}$</td>
<td>Evaporation of snow in the unsaturated sub-cloud layer</td>
</tr>
<tr>
<td>$s$</td>
<td>Freezing of condensate in the updraughts</td>
</tr>
<tr>
<td>$G_{precip}$</td>
<td>Conversion rate from cloud (water+ice) into precipitation (rain+snow)</td>
</tr>
<tr>
<td>$G_{rain}$</td>
<td>Conversion rate from cloud water into rain</td>
</tr>
<tr>
<td>$G_{snow}$</td>
<td>Conversion rate from cloud ice into snow</td>
</tr>
<tr>
<td>$h_{up}$</td>
<td>Moist static energy ($= c_pT + Lq + gz$) in the environment</td>
</tr>
<tr>
<td>$h_{sat}$</td>
<td>Saturated moist static energy in the environment</td>
</tr>
<tr>
<td>$h_{up}$</td>
<td>Moist static energy in the updraughts</td>
</tr>
<tr>
<td>$h_{down}$</td>
<td>Moist static energy in the downdraughts</td>
</tr>
<tr>
<td>$K_{up}$</td>
<td>Kinetic energy in the updraughts</td>
</tr>
<tr>
<td>$L$</td>
<td>Effective latent heat for an ice/water mix</td>
</tr>
</tbody>
</table>
### Chapter 5 ‘Convection’

- $L_{\text{fus}}$: Latent heat of fusion
- $L_{\text{subl}}$: Latent heat of sublimation
- $L_{\text{vap}}$: Latent heat of vaporization
- $l_{\text{up}}$: Cloud water/ice content in the updraughts
- $M$: Melting of snow
- $M_{\text{clu}}$: Net mass flux in the convective clouds (updraughts and downdraughts)
- $M_{\text{up}}$: Net mass flux in the updraughts
- $M_{\text{down}}$: Net mass flux in the downdraughts
- $P_{\text{rain}}$: Net flux of precipitation in the form of rain
- $P_{\text{snow}}$: Net flux of precipitation in the form of snow
- $p$: Pressure
- $q$: Specific humidity of the environment
- $q_{\text{up}}$: Weighted average specific humidity in the updraughts
- $q_{\text{down}}$: Weighted average specific humidity in the downdraughts
- $R$: Rain intensity
- $r_{\text{up}}$: Precipitation (rain+snow) in the updraughts
- $r_{\text{down}}$: Precipitation (rain+snow) in the downdraughts
- $S_{\text{fallout}}$: Fall-out of rain/snow
- $s$: Dry static energy in the environment
- $s_{\text{up}}$: Weighted average dry static energy in the updraughts
- $s_{\text{down}}$: Weighted average dry static energy in the downdraughts
- $T_{e}$: Virtual temperature in the environment
- $T_{v}$: Virtual temperature in the updraughts
- $u$: Component of wind in the environment
- $u_{\text{up}}$: Weighted average $u$ component of wind in the updraughts
- $u_{\text{down}}$: Weighted average $u$ component of wind in the downdraughts
- $V$: Mean terminal velocity of precipitation (rain+snow)
- $V_{\text{rain}}$: Mean terminal velocity of rain drops
- $v$: Component of wind in the environment
- $v_{\text{up}}$: Weighted average $v$ component of wind in the updraughts
- $v_{\text{down}}$: Weighted average $v$ component of wind in the downdraughts
- $w$: Vertical velocity in the environment
- $w_{\text{up}}$: Weighted average vertical velocity in the updraughts
- $\rho$: Density of air
- $\rho_{\text{rain}}$: Density of rain within the grid box
- $\tau$: Adjustment time scale
CHAPTER 6  Clouds and large-scale precipitation

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6.3 Code

6.1 THEORY
Cloud and large-scale precipitation processes are described by prognostic equations for cloud liquid water/ice and cloud fraction and diagnostic relations for precipitation. The scheme is described in detail in Tiedtke (1993).

6.1.1 Definitions

6.1.1 (a) Specific cloud water content and cloud fraction. The grid-mean specific cloud water/ice content is defined as

\[ l = \frac{1}{V} \int \frac{\rho_w}{\rho} dV, \]  

(6.1)

where \( \rho_w \) is the density of cloud water, \( \rho \) is the density of moist air and \( V \) is the volume of the grid box. The
fraction of the grid box covered by clouds is defined as

\[ a = \frac{1}{V} \int \delta \, dV, \quad \delta = \begin{cases} 
1, & \text{in clouds} \\
0, & \text{otherwise}
\end{cases} \quad (6.2) \]

Furthermore, the definition of the specific cloud water content per cloud area (in-cloud water/ice content) is

\[ l_{\text{cla}} = \frac{l}{a}. \quad (6.3) \]

6.1.1 (b) Saturation specific humidity. The saturation specific humidity is expressed as a function of saturation water vapour pressure as

\[ q_{\text{sat}} = \frac{R_{\text{dry}} e_{\text{sat}}(T)}{p - (1 - \frac{R_{\text{dry}}}{R_{\text{vap}}}) e_{\text{sat}}(T)}, \quad (6.4) \]

where the saturation water vapour pressure is expressed with the Tetens formula

\[ e_{\text{sat}}(T) = a_3 \exp \left( a_4 \left( \frac{T - T_0}{T - a_3} \right) \right), \quad (6.5) \]

where \( a_3 \) and \( a_4 \) are different depending on the sign of \( (T - T_0) \) (i.e. water or ice phase with \( T_0 = 273.16 \) K)

6.1.1 (c) Mixed phase. In the scheme only one variable for condensed water species is used. The distinction between the water and ice phase is made as a function of temperature. The fraction of water in the total condensate is described as

\[ \alpha = 0 \quad T \leq T_{\text{ice}}, \]
\[ \alpha = \left( \frac{T - T_{\text{ice}}}{T_0 - T_{\text{ice}}} \right)^2 \quad T_{\text{ice}} < T < T_0, \]
\[ \alpha = 1 \quad T \geq T_0. \quad (6.6) \]

\( T_{\text{sat}} \) and \( T_0 \) represent the threshold temperatures between which a mixed phase is allowed to exist and are chosen as \( T_{\text{ice}} = 250.16 \) K and \( T_0 = 273.16 \) K. The saturation thermodynamics are calculated according to the mixture of water and ice obtained with Eq. (6.6) so that the saturation specific humidity becomes

\[ q_{\text{sat}} = \alpha q_{\text{sat}(w)} + (1 - \alpha) q_{\text{sat(i)}}, \quad (6.7) \]

where \( q_{\text{sat}(w)} \) and \( q_{\text{sat(i)}} \) are the saturation specific humidities with respect to water and ice, respectively. The latent heat of phase changes is described as

\[ L = \alpha L_{\text{vap}} + (1 - \alpha) L_{\text{subl}}. \quad (6.8) \]
6.1.2 Basic equations

With these definitions and the usual assumption that clouds encountered extend vertically over the whole model layer depth the equations for the time change of the grid-box averaged cloud water/ice content and the cloud fraction are obtained as

\[
\frac{\partial l}{\partial t} = A(l) + S_{\text{conv}} + S_{\text{bl}} + S_{\text{strat}} - E_{\text{cld}} - G_{\text{prec}} - \frac{1}{\rho} \frac{\partial}{\partial z} (\rho \overline{w'\theta'})_{\text{entr}} \tag{6.9}
\]

and

\[
\frac{\partial a}{\partial t} = A(a) + \delta a_{\text{conv}} + \delta a_{\text{bl}} + \delta a_{\text{strat}} - \delta a_{\text{evap}} \tag{6.10}
\]

The terms on the right-hand side of Eq. (6.9) and Eq. (6.10) represent the following processes:

- $A(l)$, $A(a)$ — transport of cloud water/ice and cloud area through the boundaries of the grid volume
- $S_{\text{conv}}$, $\delta a_{\text{conv}}$ — formation of cloud water/ice and cloud area by convective processes
- $S_{\text{bl}}$, $\delta a_{\text{bl}}$ — formation of cloud water/ice and cloud area by boundary-layer turbulence
- $S_{\text{strat}}$, $\delta a_{\text{strat}}$ — formation of cloud water/ice and cloud area by stratiform condensation processes
- $E_{\text{cld}}$ — rate of evaporation of cloud water/ice
- $G_{\text{prec}}$ — generation of precipitation from cloud water/ice
- $(1/\rho) \partial (\rho \overline{w'\theta'})_{\text{entr}} / \partial z$ — dissipation of cloud water/ice by cloud top entrainment
- $\delta a_{\text{evap}}$ — rate of decrease of cloud area due to evaporation.

The large-scale budget equations for specific humidity $q$, and dry static energy $s = c_p T + gz$ after introduction of the scheme are modified to

\[
\frac{\partial q}{\partial t} = A(q) - S_{\text{bl}} - S_{\text{strat}} + E_{\text{cld}} + E_{\text{prec}} - \frac{1}{\rho} \frac{\partial}{\partial z} (\rho \overline{w'q'})_{\text{entr}} \tag{6.11}
\]

and

\[
\frac{\partial s}{\partial t} = A(s) + L(S_{\text{bl}} + S_{\text{strat}} - E_{\text{cld}} - E_{\text{prec}}) - L_{\text{fus}} M - \frac{1}{\rho} \frac{\partial}{\partial z} (\rho \overline{w's'})_{\text{entr}} + c_p \{(1-a) q_{\text{clear}} + a q_{\text{cld}}\} \tag{6.12}
\]

where $A(q)$ and $A(s)$ represent all processes except those related to clouds and radiation. $L_{\text{fus}}$ is the latent heat of freezing, $M$ is the rate of snowmelt, $q_{\text{clear}}$ and $q_{\text{cld}}$ are the radiative heating rates in cloud-free and cloudy areas. The flux-divergence terms represent the effects of cloud top entrainment.

6.1.3 Definition of the source and sink terms

6.1.3 (a) Convection. Clouds formed by convective processes are parametrized by considering them to be condensates produced in cumulus updraughts and detrained into the environmental air. This approach, besides being part of the cloud parametrization, represents also an important extension of the model’s cumulus parametrization. It is applied for all types of convection, namely deep, shallow and mid-level. The source of cloud water/ice content is
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\[ S_{conv} = \frac{D_{up}}{\rho} (l_{up} - l) + \frac{M_{up} \partial l}{\rho \partial z} \]  

(6.13)

and the source of cloud area is described as

\[ \partial a_{conv} = (1 - a) \frac{D_{up}}{\rho} + \frac{M_{up} \partial a}{\rho \partial z} . \]  

(6.14)

where \( D_{up} \) is the detrainment of mass from cumulus updraughts, \( l_{up} \) is the specific cloud water/ice content in cumulus updraughts and \( M_{up} \) is the updraught mass flux (see chapter 5). The factor \((1 - a)\) in Eq. (6.14) appears because updraught air detrains simultaneously into cloud-free air as well as into already existing clouds.

6.1.3 (b) Boundary layer clouds. This part of the scheme considers stratocumulus clouds at the top of convective boundary layers. They are distinguished from shallow cumuli by making the assumption, that the cloud depth must not exceed one model-layer depth. All clouds deeper than one layer are represented as convective clouds by the cumulus convection scheme. The scheme follows the mass-flux approach, so that the cloud transport for moisture is written as

\[ F_q = \rho \hat{w} (q_{up} - q_{down}) , \]  

(6.15)

where \( q_{up} \) and \( q_{down} \) are updraught and downdraught specific humidity, respectively, and \( \rho \hat{w} \) (\( = \rho a_{up} w_{up} \)) is the cloud mass flux, \( w_{up} \) being the updraught velocity and \( a_{up} \) the fractional area of updraughts. Note that in contrast to convection, stratocumulus cloud circulations contain roughly equal ascending and descending branches. The cloud-base mass flux is determined by reformulating the moisture transport at cloud base produced by the boundary layer parametrization \( J_q \) (see Sections 3.3 and 3.4 of Chapter 3 ‘Turbulent diffusion and interactions with the surface’) into the mass-flux concept so that

\[ (\rho \hat{w})_{base} = \frac{J_q}{q_0 - \{a(q_{sat} + l_{cl})_{top} + (1 - a) q_{top}\}} . \]  

(6.16)

The subscripts ‘0’ and ‘top’ refer to model levels near the surface and close to the cloud top (i.e. next level above cloud base), respectively, indicating that the updraughts start close to the surface and the downdraughts close to the cloud top. Above cloud base the assumption is made that \( \hat{w} \) decreases linearly to zero at cloud top. The net generation of cloud water/ice due to condensation in updraughts and evaporation in downdraughts then becomes

\[ S_{bl} = \frac{1}{\rho} \frac{\partial (\rho \hat{w})}{\partial z} (l_{up} - a l_{down}) , \]  

(6.17)

and the source of cloud air in terms of cloud cover is

\[ \partial a_{bl} = \frac{1}{\rho} \frac{\partial (\rho \hat{w})}{\partial z} (1 - a) . \]  

(6.18)

6.1.3 (c) Formation of stratiform clouds. Here the formation of clouds by non-convective processes (e.g. large-scale lifting of moist air, radiative cooling etc.) is considered. The parametrization is based on the principle that condensation processes are determined by the rate at which the saturation specific humidity decreases. This rate is linked to vertical motions and diabatic cooling through
\[
\frac{dq_{\text{sat}}}{dt} = \left( \frac{dq_{\text{sat}}}{dp} \right)_{\text{ma}} \left( \bar{w} + gM_{\text{Cu}} \right) + \left( \frac{dq_{\text{sat}}}{dT} \right)_{\text{diab}} \cdot
\]

where \( \left( \frac{dq_{\text{sat}}}{dp} \right)_{\text{ma}} \) is the change of \( q_{\text{sat}} \) along a moist adiabat through point \((p, T)\), \( \bar{w} \) is the area-mean generalized vertical velocity, \( gM_{\text{Cu}} \) is the cumulus-induced subsidence between the updrafts, and \( \left( \frac{dq_{\text{sat}}}{dT} \right)_{\text{diab}} \) is the net temperature tendency due to radiative and turbulent processes. Two cases of condensation are distinguished

(a) in already existing clouds

(b) the formation of new clouds

\[
c_{\text{cld}} = c_1 + c_2.
\]

Condensation in already existing clouds is described as

\[
c_1 = -\alpha \frac{dq_{\text{sat}}}{dt} \quad \frac{dq_{\text{sat}}}{dt} < 0.
\]

New clouds are assumed to form, when the grid-averaged relative humidity exceeds a threshold value which is defined as a function of height as

\[
\begin{align*}
\text{RH}_{\text{crit}} &= \text{RH}_c + (1 - \text{RH}_c) \left( \frac{\sigma - \sigma_1}{1 - \sigma_1} \right)^2 \quad \sigma_1 < \sigma \\
\text{RH}_{\text{crit}} &= \text{RH}_c \quad \sigma_{\text{trop}} + (\Delta \sigma)_d < \sigma < \sigma_d \\
\text{RH}_{\text{crit}} &= \text{RH}_c + (1 - \text{RH}_c) \left( \frac{\sigma_{\text{trop}} + (\Delta \sigma)_d - \sigma}{(\Delta \sigma)_d} \right)^2 \quad \sigma_{\text{trop}} < \sigma < \sigma_{\text{trop}} + (\Delta \sigma)_d \\
\text{RH}_{\text{crit}} &= 1 \quad \sigma < \sigma_{\text{trop}}
\end{align*}
\]

where \( \text{RH}_c = 0.8 \), \( \sigma = p/p_{\text{surf}} \) with \( p \) being the pressure and \( p_{\text{surf}} \) the pressure at the surface, \( \sigma_1 = 0.8 \), \( \sigma_{\text{trop}} \) is the height of the tropopause in \( \sigma \)-coordinates and \( (\Delta \sigma)_d = 0.2 \). The increase in cloud cover is determined by how much of the cloud-free area exceeds saturation in one time step which in turn depends on the moisture distribution in the cloud-free area and how fast saturation is approached. The moisture is assumed to be evenly distributed within the range \( \{ q_{\text{env}}(q_{\text{sat}} - q_{\text{env}}) \cdot q_{\text{sat}} \} \) around the mean environmental value \( q_{\text{env}} \), while the approach to saturation is determined by \( dq_{\text{sat}}/dt \). The increase in cloud cover then becomes

\[
\delta a_{\text{strat}} = \frac{(1 - a)}{(q_{\text{sat}} - q_{\text{env}})} \cdot \frac{dq_{\text{sat}}}{dt} \quad \frac{dq_{\text{sat}}}{dt} < 0,
\]

which can be expressed in terms of grid averages (using the definition \( q = aq_{\text{sat}} + (1 - a)q_{\text{env}} \)) as

\[
\delta a_{\text{strat}} = -(1 - a)^2 \frac{1}{(q_{\text{sat}} - q)} \cdot \frac{dq_{\text{sat}}}{dt} \quad \frac{dq_{\text{sat}}}{dt} < 0.
\]

For the application of Eq. (6.24) at values of \( q \) close to saturation, the constraint \( \delta a_{\text{strat}} < (1 - a)/\Delta t \) is imposed to ensure realistic values of \( a \).

The generation of cloud water/ice in newly formed clouds is then
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\[ c_2 = \frac{1}{2} \delta a_{\text{strat}} \frac{dq_{\text{sat}}}{dt} \quad \text{if } dq_{\text{sat}}/dt < 0 , \]  
\text{(6.25)}


where \( \delta a_{\text{strat}} \) is the fractional cloud cover produced in the time step by Eq. (6.24).

6.1.3 (d) Evaporation of cloud water/ice. The scheme describes evaporation of clouds by two processes in connection with large-scale and cumulus-induced descent and diabatic heating and by turbulent mixing of cloud air with unsaturated environmental air.

\[ \mathcal{E}_{\text{cl}} = \mathcal{E}_1 + \mathcal{E}_2. \]  
\text{(6.26)}

The first process is accounted for in the same way as stratiform cloud formation except that \( dq_{\text{sat}}/dt > 0 \). Hence

\[ \mathcal{E}_1 = a \frac{dq_{\text{sat}}}{dt} \quad \text{if } dq_{\text{sat}}/dt > 0 . \]  
\text{(6.27)}

Assuming a homogeneous horizontal distribution of liquid water in the cloud, the cloud fraction remains unaltered by this process except at the final stage of dissipation where it reduces to zero.

\[ \delta a_{\text{evap}} = \frac{a}{\Delta t} \quad \text{if } l \to 0 \]  
\text{(6.28)}

The parametrization of cloud dissipation as cloud air mixes with environmental air is described as a diffusion process proportional to the saturation deficit of the environmental air:

\[ \mathcal{E}_2 = a K (q_{\text{sat}} - q) , \]  
\text{(6.29)}

where the diffusion coefficient is

\[ K = 10^{-6} \text{ s}^{-1} . \]  
\text{(6.30)}

The decrease in cloud cover is parametrized as

\[ \delta a_{\text{evap}} = \frac{\mathcal{E}_2}{l_{\text{cl}}} , \]  
\text{(6.31)}

where \( l_{\text{cl}} \) is the specific cloud water/ice content per cloud area as defined in Eq. (6.3). Note that because of Eq. (6.3) the parametrizations Eq. (6.29) and Eq. (6.31) imply a reduction in cloud area while \( l_{\text{cl}} \) remains unchanged.

6.1.3 (e) Cloud top entrainment. Fluxes of heat, moisture, cloud water/ice, and momentum through cloud top due to the cloud top entrainment process are described as

\[ (w \Phi)'_{\text{ent}} = -w_{\text{ent}} \Delta \Phi , \]  
\text{(6.32)}

where \( \Phi \) stands for any of the transported variables and \( w_{\text{ent}} \) is the entrainment velocity. \( \Delta \) stands for the change of \( \Phi \) between two model levels. The parametrization of cloud top entrainment is currently only used if the level above a cloudy model level is entirely cloud free and if \( \Delta s_{\nu} \) is positive (stable layer), where \( s_{\nu} \) represents the vir-
tual dry static energy between the two layers. There are two parametrized contributions to the entrainment velocity

\[ w_c = w_{c,1} + w_{c,2} \]  \hfill (6.33)

(i) Clouds at the top of convective boundary layers. In the case of clouds at the top of convective boundary layers the parametrization of the entrainment velocity follows Deardorff (1976). The entrainment velocity is represented as

\[ w_{c,1} = \frac{-(\overline{w's_v})_H}{\Delta s_v} \]  \hfill (6.34)

where

\[ -(\overline{w's_v})_H = \alpha \frac{1}{H} \int_{0}^{H} \overline{w's_v} \, dz \]  \hfill (6.35)

is the average buoyancy flux in the mixed layer of height \( H \) and \( \alpha = 0.5 \).

(ii) All cloud tops. The second contribution to the entrainment velocity is parametrized as

\[ w_{c,2} = \beta \frac{c_{LW}}{\Delta s_v} \]  \hfill (6.36)

where \( \Delta f_{LW} \) is the longwave radiative flux divergence and \( \beta = 0.5 \).

Cloud water/ice transported into the cloud free layer above by entrainment is assumed to evaporate immediately.

6.1.3 (f) Precipitation processes. Similar to radiation, precipitation processes are treated separately in clear and cloudy skies. This owes to the fact that the microphysical processes in these two regions are very distinct from each other, with conversion, collection and accretion processes being relevant in clouds whereas evaporation of precipitation is the relevant process outside clouds. Therefore the precipitation flux is written as

\[ P = P^{clld} + P^{clr} \]  \hfill (6.37)

with

\[ P^{clld} = \frac{1}{A} \int P \cdot H(l) \, dA \]  \hfill (6.38)

and

\[ P^{clr} = \frac{1}{A} \int P \cdot (1 - H(l)) \, dA \]  \hfill (6.39)

where the step function, \( H(l) \), marks the portion of the grid-cell containing cloud with a condensate specific humidity \( l > 0 \) and \( A \) is the area of the grid-cell.

The precipitation fraction in the gridbox is then described as
Precipitation sources are represented differently for pure ice clouds and for mixed phase and pure water clouds.

\[ G_{\text{prec}} = G_t + G_{mw} \]  \hspace{1cm} (6.43)

The distinction is made as a function of temperature according to Eq. (6.6). The rain and snow formed is removed from the column immediately but can evaporate, melt and interact with the cloud water in the layers it passes through.

(i) \textit{Pure ice clouds.} The precipitation process in ice clouds is treated separately for two classes of particles. The separation is made by size at a threshold of 100 \( \mu \text{m} \). First the ice water content in particles smaller than 100 \( \mu \text{m} \) is determined following a parametrization proposed by McFarquhar and Heymsfield (1997) as

\[ IWC_{c100} = \frac{1}{1000} \cdot \min \left[IWC_{\text{tot}}, b_1 \left( \frac{IWC_{\text{tot}}}{IWC_0} \right)^{b_2} \right] \]  \hspace{1cm} (6.44)

where

\[ IWC_{\text{tot}} = \frac{\rho_{l,cld}}{1000} \]  \hspace{1cm} (6.45)

is the total ice water content in \( \text{g m}^{-3} \). \( IWC_0 \) is set to 1 \( \text{g m}^{-3} \), \( b_1 = 0.252 \text{ g m}^{-3} \) and \( b_2 = 0.837 \). The fallout of the so diagnosed \( IWC_{c100} \) (now in \( \text{kg m}^{-3} \)) is treated as a sedimentation of the ice particles with a terminal fall speed of

\[ w_{\text{ice}} = c_1 IWC_{c100}^{c_2} \]  \hspace{1cm} (6.46)

based on Heymsfield and Donner (1990); the constants currently chosen are \( c_1 = 3.29 \) and \( c_2 = 0.16 \). The ice content in particles larger than 100 \( \mu \text{m} \)

\[ IWC_{>100} = 1000 \cdot IWC_{\text{tot}} - IWC_{c100} \]  \hspace{1cm} (6.47)

is converted into snow within one timestep.
Given the fallspeed and the separation by particle size the contribution to $G_{\text{prec}}$ from pure ice clouds can be written as

$$G_i = \frac{1}{\rho} \frac{\partial}{\partial z} (w_{ic} \text{IWC}_{<100}) + \frac{\text{IWC}_{>100}}{\tau},$$  \hspace{1cm} (6.48)

where $\tau$ is the time-step of the model. For the small particle ice settling into cloudy area is treated as source of cloud ice in the layer below whereas ice settling into clear sky is converted into snow (see Subsection 6.2.3). Note that the minus sign in the first term of the right hand side of (6.48) appears since the fall velocity of ice is assumed to be positive downwards.

(ii) Mixed phase and pure water clouds. For mixed phase and pure water clouds a parametrization following Sundqvist (1978) is used. The generation of precipitation is written as

$$G_{mw} = a c_0 l_{\text{clad}} \left[ 1 - \exp \left\{ -\left( \frac{l_{\text{clad}}}{l_{\text{crit}}^*} \right)^2 \right\} \right],$$  \hspace{1cm} (6.49)

where $c_0^{-1}$ represents a characteristic time scale for conversion of cloud droplets into drops and $l_{\text{crit}}$ is a typical cloud water content at which the release of precipitation begins to be efficient. These disposable parameters are adjusted as follows

$$c_0 = c_0^* F_1 F_2$$  \hspace{1cm} (6.50)

and

$$l_{\text{crit}} = \frac{l_{\text{crit}}^*}{F_1 F_2}$$  \hspace{1cm} (6.51)

to take into account the effect of collection of cloud droplets by raindrops falling through the cloud $(F_1)$ and the Bergeron–Findeisen mechanism $(F_2)$. Here $F_1$ and $F_2$ are defined as

$$F_1 = 1 + b_1 \sqrt{P_{\text{loc}}},$$  \hspace{1cm} (6.52)

and

$$F_2 = 1 + b_2 \sqrt{(T_{\text{BF}} - T)} \quad \text{if} \quad T_{\text{ice}} < T < T_{\text{BF}}$$  \hspace{1cm} (6.53)

where $P_{\text{loc}}$ is the local cloudy precipitation rate $(P_{\text{loc}} = P_{\text{clad}}^d / a_p^{clad})$ and $T_{\text{BF}}$ is the temperature at which the Bergeron–Findeisen mechanism starts to enhance the precipitation. The values for the constants are those used by Sundqvist (1978), namely $T_{\text{BF}} = 268 \text{ K}$, $b_1 = 100$, $b_2 = 0.5$, $c_0^* = 10^{-3} \text{ s}^{-1}$, and $l_{\text{crit}}^* = 0.3 \text{ g kg}^{-1}$.

(iii) Evaporation of precipitation. The parametrization of rain and snow evaporation is uncertain. A scheme following Kessler (1969) is used. It describes the evaporation rate as
where \( a_p^{clr} \) is the clear-sky precipitation fraction. Evaporation of rain/snow only takes place when the grid mean relative humidity is below a threshold value. The choice of the threshold value is not straightforward for numerical reasons. Here, the assumption is made that the clear-sky relative humidity (= grid mean relative humidity in the absence of clouds) that can be reached by evaporation of precipitation is a function of the fractional coverage with precipitation of the clear sky part of the grid-box. Hence, the threshold value is parametrized as

\[
RH_{crit,p} = 0.7 + 0.3 \frac{a_p^{clr} \frac{1}{1 - a}}{1 - a}.
\]  

(iv) Melting of snow. The melting of snow is parametrized by allowing the part of the grid box that contains precipitation to cool to \( T_{melt} \) over a time scale \( \tau \), i.e.,

\[
\mathcal{M} = (a_p^{cl} + a_p^{clr}) \frac{c_p}{L_{ fus}} \frac{T - T_{melt}}{\tau},
\]

where \( T_{melt} = 0^\circ C \) and

\[
\tau = \frac{5h}{1 + 0.5(T - T_{melt})}.
\]

6.2 Numerics

6.2.1 Integration of the equations for cloud water/ice and cloud cover

As cloud processes are rapidly varying in time, care must be taken when Eq. (6.9) and Eq. (6.10) are integrated over the relatively large model time steps. Therefore terms that depend linearly on \( a \) and \( l \) are integrated analytically. Eq. (6.9) and Eq. (6.10) can be written as

\[
\frac{\partial l}{\partial t} = C - Dl
\]  

where C is defined by Eq. (6.13), Eq. (6.17), Eq. (6.21), Eq. (6.25), Eq. (6.27) and Eq. (6.29) and D is defined by Eq. (6.48) or Eq. (6.49) respectively, and

\[
\frac{\partial a}{\partial t} = (1 - a)A - aB + C_a
\]

with \( A \), \( B \), and \( C_a \) defined by Eq. (6.14), Eq. (6.18), Eq. (6.24), Eq. (6.28) and Eq. (6.31). Analytical integration of Eq. (6.57) and Eq. (6.58) yields

\[
a(t + \Delta t) = a(t) \exp \{- (A + B) \Delta t \} + \frac{A + C_a}{A + B} [1 - \exp \{- (A + B) \Delta t \}]
\]
and

\[ l(t + \Delta t) = l(t) \exp(-D\Delta t) + \frac{C}{D} \{ 1 - \exp(-D\Delta t) \} \]  

(6.60)

Since there are terms in Eq. (6.57) that depend on \(a\) and terms in Eq. (6.58) that depend on \(l\) the following method to solve the two equations is adopted. First Eq. (6.59) is solved using the values for \(l\) at time \(t\). Then a time-centred value for the cloud fraction, \(\hat{a}\) is calculated as

\[ \hat{a} = \frac{a(t + \Delta t) + a(t)}{2}. \]  

(6.61)

Then Eq. (6.60) is divided by \(\hat{a}\) and solved yielding a new value for the in-cloud value \(l_{\text{cl}}(t + \Delta t)\) which is converted into the grid-mean value by re-multiplying with \(\hat{a}\).

### 6.2.2 Calculation of \(dq_{\text{sat}}/dt\)

Special care has to be taken in the numerical calculation of \(dq_{\text{sat}}/dt\) from Eq. (6.19). Since the saturation water vapour pressure depends exponentially on temperature, straightforward numerical integration of Eq. (6.19) would produce large truncation errors. Therefore the average of \(dq_{\text{sat}}/dt\) over the time step is determined by the means of moist adjustment (e.g. Halldin and Williams 1980). This is achieved by first extrapolating the cloud temperature to time-level \(t + \Delta t\) and then adjust temperature and moisture toward saturation conditions.

### 6.2.3 Convective cloud source

The vertical discretization of equations (6.13) and (6.14) is achieved with a simple upstream scheme, i.e.,

\[ S_{\text{conv}} = \frac{D_{up}}{p_k} (l_{up,k+1/2} - l_k) - g M_{up,k-1/2} \frac{l_{k-1} - l_k}{p_{k-1} - p_k} \]  

(6.62)

and

\[ \delta a_{\text{conv}} = \frac{D_{up}}{p_k} (1 - a_k) - g M_{up,k-1/2} \frac{a_{k-1} - a_k}{p_{k-1} - p_k}. \]  

(6.63)

Although two of the terms in equation (6.62) depend linearly on \(l_k\) it was decided to treat the convective source (like any other source of condensate) fully explicitly, i.e., (6.62) is added into (6.57) as a contribution to \(C_k\) only.

For cloud fraction it is obvious that the first term on the right hand side of (6.63) can be added to \(A_k\) in equation (6.58) whereas the second term can be split into a contribution to \(B_k\) and \(C_{a,k}\).

### 6.2.4 Stratiform cloud source

It is evident from (6.24) that the stratiform source of cloud cover is quadratically dependent on \((1 - a)\) and can therefore not easily be integrated analytically following (6.58). To overcome this problem one factor of \((1 - a)\) is integrated into \(A_k\), i.e., treated explicitly, before carrying out the analytic integration of (6.58), i.e.,

\[ A_{k,\text{strat}} = \frac{1 - a_k}{(q_{\text{sat},k} - q_k)} \left( \frac{dq_{\text{sat}}}{dt} \right)_k. \]  

(6.64)
6.2.5 Precipitation fractions

The method to determine \( a_P^{cl} \) and \( a_P^{clr} \) is as follows. If precipitation is generated in a level through the processes of autoconversion or ice sedimentation, it is assumed to be generated at all portions of the cloud uniformly and thus at the base of level \( k, a_P^{cl} = a_k \). The precipitation generated in this cloudy region is given by:

\[
\Delta P_k^{cl} = \frac{1}{\Delta A} \left( \int_0^{P_{k+1/2}} G_{prec} \cdot H(l) dp \right) dA ,
\]

and the cloudy precipitation flux at the base of level \( k \) is given by \( P_k^{cl} = \tilde{P}_k^{cl} + \Delta P_k^{cl} \), where the twiddle symbol indicates the value of \( P_k^{cl} \) at the top of level \( k \). Because the cloud is assumed to be internally homogenous, (6.65) simplifies to \( a_k G_{prec}(p_{k+1/2} - p_{k-1/2})/g \), where \( \tilde{S}_p^{cl} \) is the generation rate of precipitation inside the cloud. If only accretion occurs in the clouds of level \( k, a_P^{cl} = \tilde{a}_P \), the fractional area that contains cloudy precipitation flux at the top of level \( k \).

Because the clear precipitation flux is assumed to be horizontally uniform, evaporation does not alter the area containing clear precipitation flux such that \( a_P^{cl} = a_k \). Only in the case that all of the clear precipitation flux evaporates in level \( k \) does \( a_P^{clr} = 0 \). The clear-sky precipitation flux at the base of level \( k \) is given by \( P_k^{clr} = \tilde{P}_k^{clr} + \Delta P_k^{clr} \), where \( \tilde{P}_k^{clr} \) is the clear-sky precipitation flux at the top of level \( k \), and

\[
\Delta P_k^{clr} = \frac{1}{\Delta A} \left( \int_0^{P_{k+1/2}} E_{prec} \cdot (1 - H(l)) dp \right) dA = \tilde{a}_P^{clr} E_{prec}(p_{k+1/2} - p_{k-1/2})/g ,
\]

where \( E_{prec} \) represents precipitation evaporation. Note that precipitation evaporation is a function of \( \tilde{P}_k \), guaranteeing that precipitation generated in a level cannot evaporate in the same level. This will guarantee consistency with the assumption that clouds where present fill the vertical extent of the grid cell and that horizontal transfer of precipitation mass from cloudy to clear regions of the grid cell is not possible.

At the interfaces between levels, precipitation mass that is in cloud of the upper level may fall into clear air of the lower level, or precipitation mass that is in clear air of the upper level may fall into cloud of the lower level. Thus at level interfaces an algorithm is needed to transfer precipitation and its area between the cloudy and clear portions of the grid box. The algorithm is constructed by determining the amount of area associated with each transfer and then transferring precipitation fluxes between clear and cloudy components according to the assumption that the precipitation flux is horizontally uniform but with different values in the clear and cloudy regions containing precipitation.

There are four possible areas to be defined (see schematic in Figure 20): the area in which cloudy precipitation flux falls into cloud of the lower level, the area in which cloudy precipitation flux falls into clear air of the lower level, the area in which clear precipitation flux falls into clear air of the lower level, and the area in which clear precipitation flux falls into cloud of the lower level. To determine these areas, the cloud overlap assumption is applied to determine the relative horizontal placements of clouds in the upper and lower levels. For the ECMWF model, the cloud overlap assumption is expressed in terms of an equation which relates the total horizontal area \( C \) covered by clouds in levels 1 to \( k \) (where \( k = 1 \) is the top level of the model), to the total horizontal area cover by clouds in levels 1 to \( k-1 \):

\[
(1 - C_k) = (1 - C_{k-1}) \cdot \frac{1 - \max(a_k, a_{k-1})}{1 - \min(a_k, a_{k-1})}.
\]
where $\delta$ is a tiny number set to $10^{-6}$. Equation (6.67) gives maximum overlap for clouds in adjacent levels and random overlap for clouds separated by clear levels. From this equation, one can determine the portion of clouds of the lower level which is unoverlapped by clouds at all higher levels; this area, $\Delta C = C_k - C_{k-1}$, cannot have any precipitation falling into it. Using this assumption, the area for which cloudy precipitation flux falls into clear air of the level below is given by:

$$\Delta a_{cld\rightarrow clr} = a^{cld}_{P,k-1} - \min(a_k - \Delta C, a^{cld}_{P,k-1}).$$

Equation (6.68) makes the further assumption that there is maximum overlap between the area covered by cloudy precipitation at the base of the upper level and the portion of the lower level cloud which lies beneath clouds in higher levels, $a_k - \Delta C$. With the assumption that the precipitation flux is horizontally uniform, the amount of cloudy precipitation flux of the upper level that falls into clear air of the level below is:

$$\Delta P_{cld\rightarrow clr} = \frac{\Delta a_{cld\rightarrow clr}}{a^{cld}_{P,k-1}} \cdot P^{cld}_{k-1}.\tag{6.69}$$

The area in which clear precipitation flux of the upper level falls into cloud of the level below is:

$$\Delta a_{clr\rightarrow cld} = \max(0, \min(a^{clr}_{P,k-1}, a_k - \Delta C - a_{k-1})),\tag{6.70}$$

which assumes maximum overlap between the portion of the cloud in the lower level $k$ which has cloud at some higher level other than $k-1$, and the area covered by the clear precipitation flux. Again, with the assumption that the precipitation flux is horizontally uniform, the amount of clear precipitation flux of the upper level that falls into cloud of the level below is:

$$\Delta P_{clr\rightarrow cld} = \frac{\Delta a_{clr\rightarrow cld}}{a^{clr}_{P,k-1}} \cdot P^{clr}_{k-1}.\tag{6.71}$$

Finally, the areas and fluxes at the top of level $k$ can be related to those at the base of level $k-1$ by:

$$\tilde{a}^{cld}_{P,k} = a^{cld}_{P,k-1} + \Delta a_{clr\rightarrow cld} - \Delta a_{cld\rightarrow clr},\tag{6.72}$$

$$\tilde{a}^{clr}_{P,k} = a^{clr}_{P,k-1} - \Delta a_{cld\rightarrow cld} + \Delta a_{clr\rightarrow cld},\tag{6.73}$$

$$\tilde{P}^{cld}_k = P^{cld}_{k-1} + \Delta P_{clr\rightarrow cld} - \Delta P_{cld\rightarrow clr},\tag{6.74}$$

$$\tilde{P}^{clr}_k = P^{clr}_{k-1} - \Delta P_{cld\rightarrow clr} + \Delta P_{clr\rightarrow cld}.$$\tag{6.75}

From these equations it is obvious that total precipitation area, $a^{cld}_{P} + a^{clr}_{P}$, and precipitation flux, $P^{cld} + P^{clr}$, are conserved at level interfaces.

### 6.2.6 Precipitation sources

After the integration of Eq. (6.60) the fallout of condensate (represented by the term $Dl$ in Eq. (6.57)) out of model
level \( k \) is determined as

\[
G_{\text{fallout},k} = \frac{l_k(t) - l_k(t + \Delta t)}{\Delta t} + C_k.
\]  

(6.76)

The condensate falling out of model level \( k \) is then distributed into rain, snow or cloud ice in the level below using the following assumptions:

1. **Pure water clouds.** In the case of pure water clouds (\( T_k \geq T_0 \)) all condensate falling out of a model level is converted into rain, i.e.,

\[
G_{\text{rain},k} = G_{\text{fallout},k}.
\]  

(6.77)

2. **Mixed phase clouds.** In the case of mixed phase clouds (\( T_{\text{ice}} < T_k < T_0 \)) all condensate falling out of a model level is converted into rain or snow whereby the partitioning between the two phases is determined using Eq. (6.6), i.e.,

\[
G_{\text{rain},k} = \alpha_k G_{\text{fallout},k}
\]  

(6.78)

and

\[
G_{\text{snow},k} = (1 - \alpha_k)G_{\text{fallout},k}.
\]  

(6.79)

3. **Pure ice clouds.** In the case of pure ice clouds (\( T_k \leq T_{\text{ice}} \)) the condensate falling out of a model level is partitioned into a source of cloud ice in the level below and snow. There are two sources of snow from falling cloud ice; i) all ice content in particles larger than 100 \( \mu m \) is converted into snow, and ii) of the falling cloud ice in particles smaller than 100 \( \mu m \), ice falling into clear sky is converted into snow, while ice in falling into cloud remains cloud ice. This is implemented in the code as follows. First (6.44) is solved to determine the ice water content in particles smaller than 100 \( \mu m \). Then (6.60) is solved for layer \( k \) using

\[
C_k = C_{\text{noprec},k} + C_{\text{fallin},k}
\]  

(6.80)

and

\[
D_k = -\alpha_{\text{<100}} \left( \frac{p_k w_{\text{ice},k}}{p_k^1 - p_{k+1}^1} + (1 - \alpha_{\text{<100}}) \frac{1}{\Delta t} \right)
\]  

(6.81)

where

\[
\alpha_{\text{<100},k} = \frac{IWC_{\text{<100},k}}{IWC_{\text{tot},k}}
\]  

(6.82)

\( C_{\text{noprec},k} \) represents all sources and sinks of cloud ice not related to precipitation processes, whereas \( C_{\text{fallin},k} \) is the source of cloud ice through settling of particles from the layer above (for definition see below). Then (6.76) is solved to determine the fallout of ice out of layer \( k \), \( G_{\text{fallout},k} \). Of the ice water content falling into layer \( k \) in particles smaller than 100 \( \mu m \), i.e.,
\( \alpha_{\text{fall},k}G_{\text{fallout},k-1} \), the part falling into overlapping cloud area is treated as source of cloud ice, \( C_{\text{fallin},k} \). The area of cloud overlap is determined as

\[
a_{\text{ovlp},k} = \min(a_k - \Delta C, a_{k-1}) .
\]

(6.83)

where \( \Delta C \) is the change of total cloud cover from layer \( k-1 \) to layer \( k \) as described above. Hence,

\[
C_{\text{fallin},k} = a_{\text{ovlp}}\alpha_{\text{fall},k}G_{\text{fallout},k-1} .
\]

(6.84)

With these definitions the generation of snow in level \( k \) becomes

\[
G_{\text{snow},k} = (1 - a_{\text{ovlp},k+1}a_{\text{fall},k})G_{\text{fallout},k} .
\]

(6.85)

After the definitions above the precipitation at the surface can be written as

\[
P_{\text{sfc}} = \int z(G_{\text{rain}} + G_{\text{snow}} - p_{\text{prec}}) \rho dz .
\]

(6.86)

6.2.7 Evaporation of precipitation

Since the evaporation of precipitation has a threshold value of relative humidity at which the process should cease to exist (see question (6.55)) an implicit treatment is applied when solving (6.54). If (6.54) is written as

\[
\frac{\partial q}{\partial t} = \beta(q_s - q) .
\]

(6.87)

the implicit solution becomes

\[
\Delta q = \frac{\beta \Delta t(q_s^n - q^n)}{1 + \beta \Delta t \left[ 1 + \frac{L}{c_p \frac{dT}{dt}} \right]^{n/\alpha(T)}} ,
\]

(6.88)

where \( n \) refers to the time level at the beginning of timestep \( \Delta t \). (6.88) ensures that evaporation of precipitation never leads to \( q_s^{n+1} > q_s^n \). To ensure the maximum relative humidity after evaporation does not exceed the threshold value defined in (6.55) the maximum change in specific humidity is calculated as

\[
(\Delta q)_{\text{max}} = \frac{RH_{\text{crit},E} \cdot q_s^n - q^n}{1 + RH_{\text{crit},E} \cdot \frac{L}{c_p \frac{dT}{dt}}} .
\]

(6.89)

The smaller of the values given by (6.88) and (6.89) is then chosen as the true value of evaporation of precipitation.

6.2.8 Cloud top entrainment

After parametrizing the entrainment flux as in Eq. (6.32) and the entrainment velocity as in Eq. (6.33) the tendency
equations for the two levels involved in the entrainment process are solved simultaneously using an implicit formulation.

The tendency equation for the cloudy model level, $k$, can be written as

$$\left( \frac{\partial \Phi}{\partial t} \right)_k = g \frac{F_{\Phi, \frac{k-1}{2}} - F_{\Phi, \frac{k+1}{2}}}{p_{\frac{k+1}{2}} - p_{\frac{k-1}{2}}}, \quad (6.90)$$

where $F_{\Phi, \frac{k-1}{2}}$ is the flux of $\Phi$ taken at half-level $k - 1/2$. A similar equation can be written for the level immediately above the cloud, $k - 1$. Since only the transport between levels $k$ and $k - 1$ are considered only the flux at half level $k - 1/2$ is non-zero. The solution for $\Phi$ at both model levels for time $n + 1$ given the values at time $n$ can then be found by solving the system of two linear equations

$$\Phi_k^{n+1} - \Phi_k^n = -g\Delta t \frac{w_e(\Phi_{k-1}^{n+1} - \Phi_k^{n+1})}{p_{\frac{k-1}{2}} - p_{\frac{k+1}{2}}} \quad (6.91)$$

and

$$\Phi_{k-1}^{n+1} - \Phi_{k-1}^n = g\Delta t \frac{w_e(\Phi_{k}^{n+1} - \Phi_{k-1}^{n+1})}{p_{\frac{k+1}{2}} - p_{\frac{k-1}{2}}} \quad (6.92).$$

### 6.2.9 Final moist adjustment

After the calculation of the liquid water/ice tendency and the corresponding tendencies of temperature and moisture a final test for supersaturation is performed. If any supersaturation is found the grid box is re-adjusted to saturation (using the moist adjustment formulation) and the moisture excess is converted into precipitation. With the introduction of the SLAVEPP method the supersaturation adjustment is applied on the effective profiles of temperature and humidity after all physical and explicit dynamical contributions have been computed. See Section 3.10 of Chapter 3 “Semi-Lagrangian formulation” for a more detailed description of the applied saturation check. Note, that small amounts of non-physical supersaturation due to numerical approximations may remain in the postprocessed fields.

### 6.3 CODE

The parametrization of cloud and large-scale precipitation processes is performed in the following routines:

**CLDPP**

This routine prepares the cloud variables for radiation calculations and calculates total, high, mid-level and low cloud cover for postprocessing.

**CLOUDSC**

This routine carries out all calculations necessary to solve Eq. (6.9) and Eq. (6.10). The calculations are carried out in the following order:

* initial setup including calculation of
  * $q_{sat}$
- tropopause height for Eq. (6.22)
- mixed layer buoyancy integral as defined in Eq. (6.35)
- convective source terms including freezing if different mixed phase assumptions are used for convection and large-scale processes (Eq. (6.13) and Eq. (6.14))
- entrainment velocity due to longwave cooling (Eq. (6.36))
- generation of clouds at top of convective boundary layer (Eq. (6.17) and Eq. (6.18))
- erosion of clouds by turbulent mixing (Eq. (6.29) and Eq. (6.31))
- calculation of $\frac{dq_{sat}}{dt}$ (see section 6.2.2)
- large-scale evaporation (Eq. (6.27))
- large-scale cloud formation (Eq. (6.21), Eq. (6.24), and Eq. (6.25))
- analytical integration of the equation for $a$ (Eq. (6.59))
- moist adjustment (see section 6.2.9)
- analytical integration of the equation for $l$ (Eq. (6.60)) and precipitation (Eq. (6.37) to Eq. (6.56) and Eq. (6.65) to Eq. (6.86)).
- melting of snow (Eq. (6.56))
- evaporation of precipitation (Eq. (6.54))
- final tendency calculations
- mixing due to cloud top entrainment (Eq. (6.91) and Eq. (6.92))
- flux calculations for diagnostics

APPENDIX A LIST OF SYMBOLS

- $A(\ )$: advective transport through the boundaries of the grid box
- $a$: fraction of grid box covered by clouds
- $a_{prec}$: fraction of grid box covered by precipitation
- $a_{up}$: fractional area of updraughts
- $c_{cl}$: condensation rate
- $c_{p}$: specific heat at constant pressure
- $D_{up}$: detrainment in the cumulus updraughts
- $\xi_{cl}$: rate of evaporation of cloud water/ice
- $\xi_{prec}$: rate of evaporation of precipitation
- $e_{sat}$: saturation water vapour
- $f_{LW}$: longwave radiative flux divergence
- $F_{q}$: moisture transport by clouds
- $g$: acceleration of gravity
- $G_{fallout}$: generation of precipitation that falls out from one level to another
- $G_{prec}$: generation of precipitation from cloud water/ice
- $G_{rain}$: generation of precipitation in the form of rain
- $G_{snow}$: generation of precipitation in the form of snow
- $H$: mixed-layer height
- $J_{q}$: surface humidity flux
- $K$: diffusion coefficient
- $L$: latent heat
- $L_{fus}$: latent heat of fusion
- $L_{subl}$: latent heat of sublimation
- $L_{vap}$: latent heat of vaporization
Part IV: ‘Physical processes’

\[ l \] grid-mean specific cloud liquid-water and ice content
\[ I_{\text{clu}} \] specific cloud water content per cloud area
\[ I_{\text{down}} \] specific cloud water/ice content in the cumulus downdraughts
\[ I_{\text{up}} \] specific cloud water/ice content in the cumulus updraughts
\[ \delta \] rate of snowmelt
\[ M_{\text{Cu}} \] cumulus-induced subsidence mass flux
\[ P \] precipitation rate
\[ P_{\text{loc}} \] local precipitation rate
\[ p \] pressure
\[ q_{\text{env}} \] environmental specific humidity
\[ q_{\text{down}} \] specific humidity in the convective downdraughts
\[ q_{\text{sat}} \] saturation specific humidity
\[ q_{\text{sat(i)}} \] saturation specific humidity with respect to ice
\[ q_{\text{sat(w)}} \] saturation specific humidity with respect to water
\[ q_{\text{up}} \] specific humidity in the convective updraughts
\[ R_{\text{clu}} \] radiation heating rate in cloudy air
\[ R_{\text{clear}} \] radiation heating rate in cloud-free air
\[ R_{\text{dry}} \] gas constant for dry air
\[ R_{\text{vap}} \] gas constant for water vapour
\[ \text{RH}_c \] = 0.8
\[ \text{RH}_{\text{crit}} \] threshold value of the relative humidity
\[ S_{\text{conv}} \] formation of cloud water/ice by convective processes
\[ S_{\text{strat}} \] formation of cloud water/ice by stratiform condensation processes
\[ S_{\text{bl}} \] formation of cloud water/ice by boundary-layer processes
\[ s \] dry static energy
\[ s_v \] virtual dry static energy
\[ T \] temperature
\[ T_0 \] = 273.16 K
\[ T_{\text{RF}} \] = 268 K temperature at which the Bergeron–Findeison enhances the precipitation
\[ T_{\text{ice}} \] = 250.16 K
\[ T_{\text{melt}} \] = 0°C
\[ \bar{w} \] area-mean generalized vertical velocity
\[ w \] \( \rho w = \rho a_{\text{up}} u_{\text{up}} \) is the cloud mass flux
\[ w_e \] entrainment velocity
\[ w_{\text{ice}} \] terminal fall speed of ice particles
\[ u_{\text{up}} \] updraught velocity
\[ \alpha \] fraction of condensate held as liquid water
\[ \delta a_{\text{bl}} \] rate of increase of cloud area by boundary-layer processes
\[ \delta a_{\text{conv}} \] rate of increase of cloud area by convective processes
\[ \delta a_{\text{strat}} \] rate of increase of cloud area by stratiform condensation processes
\[ \delta a_{\text{evap}} \] rate of decrease of cloud area due to evaporation
\[ \rho \] density of moist air
\[ \rho_w \] density of cloud water
CHAPTER 7  Surface parametrization

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7.1 INTRODUCTION

The parametrization scheme described in this chapter represents the surface fluxes of energy and water and, where appropriate, corresponding sub-surface quantities. Fig. 7.1 summarizes the main features of the land part of the model; hereafter the scheme will be referred to as the TESSEL (Tiled ECMWF Scheme for Surface Exchanges over Land) scheme. At the interface between the surface and the atmosphere, each grid-box is divided into fractions (tiles), with up to 6 fractions over land (bare ground, low and high vegetation, intercepted water, shaded and ex-
posed snow) and up to 2 fractions over sea and freshwater bodies (open and frozen water). Each fraction has its own properties defining separate heat and water fluxes used in an energy balance equation solved for the tile skin temperature. Special attention is devoted to the different physical mechanisms limiting evaporation of bare ground and vegetated surfaces.

![Schematics of the land surface](image)

Figure 7.1 Schematic representation of the structure of TESSEL land-surface scheme

Over land, the skin temperature is in thermal contact with a four-layer soil or, if there is snow present, a single layersnow mantle overlying the soil. The snow temperature varies due to the combined effect of top energy fluxes, basal heat flux and the melt energy. The soil heat budget follows a Fourier diffusion law, modified to take into account the thermal effects of soil water phase changes. The energy equation is solved with a net ground heat flux as the top boundary condition and a zero-flux at the bottom.

Snowfall is collected in the snow mantle, which in turn is depleted by snowmelt, contributing to surface runoff and soil infiltration, and evaporation. A fraction of the rainfall is collected by an interception layer, where the remaining fraction (throughfall) is partitioned between surface runoff and infiltration. Subsurface water fluxes are determined by Darcy’s law, used in a soil water equation solved with a four-layer discretization shared with the heat budget equation. Top boundary condition is infiltration plus surface evaporation, free drainage is assumed at the bottom; each layer has an additional sink of water in the form of root extraction over vegetated areas.

Finally, open water points have a fixed surface temperature. When present, frozen water occupies a fraction of the
grid box, with a prognostic ice temperature evolving in the forecast following the heat budget of a four-layer ice model in thermal contact with an underlying ocean at freezing temperature.

7.2 TILES AND SURFACE FLUXES

7.2.1 Tile and vegetation characteristics

Grid-box surface fluxes are calculated separately for the different subgrid surface fractions (or “tiles”), leading to a separate solution of the surface energy balance equation and skin temperature for each of these tiles. This is an analogue of the “mosaic” approach of Koster and Suarez (1992). Note that the tiles at the interface soil–atmosphere are in energy and hydrological contact with one single atmospheric profile above and one single soil profile below. Each grid box is divided into 8 fractions: two vegetated fractions (high and low vegetation without snow), one bare soil fraction, three snow/ice fractions (snow on bare ground/low vegetation, high vegetation with snow beneath, and sea-ice, respectively), and two water fractions (interception reservoir, ocean/lakes). The tile for “high vegetation with snow beneath” is a combined tile with a separate energy balance and evaporation model for the high vegetation and the underlying snow. A mixture of land and water (ocean/inland water) tiles is not allowed, i.e. a grid box is either 100% land or 100% sea.

In each grid box two vegetation types are present: a high and a low vegetation type. An external climate database, based on the Global Land Cover Characteristics (GLCC) data that has been derived using one year of Advanced Very High Resolution Radiometer (AVHRR) data and ancillary information (Loveland et al. 2000; http://edcdaac.usgs.gov/glcc/glcc.html; see also Chapter 10). The nominal resolution is 1 km. The data used provides for each pixel a biome classification based on the Biosphere–Atmosphere Transfer Scheme (BATS) model (Dickinson et al. 1993), and four parameters have been derived for each grid box: dominant vegetation type, \( T_H \) and \( T_L \), and the area fraction, \( A_H \) and \( A_L \), for each of the high- and low-vegetation components, respectively.

The coverage \( C_i \) for the tile \( i \) depends on the type and relative area of low and high vegetation, and the presence of snow and intercepted water. In the absence of snow and interception, the vegetation coverage of high (\( c_H \)) and low (\( c_L \)) vegetation are calculated as \( A_H c_{\text{veg}}(T_H) \) and \( A_L c_{\text{veg}}(T_L) \), respectively, with \( c_{\text{veg}} \) a vegetation type dependent coverage (see Table 7.1). The bare ground fraction \( c_B \) is the residual.

\[
\begin{align*}
    c_H &= A_H c_{\text{veg}}(T_H) \\
    c_L &= A_L c_{\text{veg}}(T_L) \\
    c_B &= 1 - c_H - c_L
\end{align*}
\] (7.1)

Each vegetation type is characterized by a series of (fixed) parameters as detailed in Table 7.1:

- A minimum canopy resistance, \( r_{c,\text{min}} \);
- A leaf area index, \( \text{LAI} \);
- A vegetation coverage, \( c_{\text{veg}} \);
- A coefficient, \( g_b \), for the dependence of the canopy resistance, \( r_c \), on water vapour pressure deficit;
- The root distribution over the soil layers, specified by an exponential profile involving attenuation coefficients, \( a_r \) and \( b_r \).

The numerical values for the parameters of Table 1 are based both on experiments conducted as described in van den Hurk et al. (2000) and on literature review, in particular Mahfouf et al. (1995), Manzi and Planton (1994), Girard and Bazile (2000), Dorman and Sellers (1989), Bonan (1994), Pitman et al. (1991), and Zeng et al. (1998).
The presence of snow and intercepted water dynamically modifies the coverage fractions. The coverage of snow, \( e_{\text{sn}} \), is linearly related to the snow mass per unit area (abbreviated to snow mass in the following), \( S \) (units \( 10^3 \text{ kg m}^{-2} \) or m). The interception reservoir fraction, \( c_i \), is given by \( W_i / W_{\text{m}} \), with \( W_{\text{m}} \), the maximum value for the intercepted water in the grid box, defined from the leaf area index contributions from the high and low vegetation tiles. The water contents of the interception reservoir, \( W_i \) (units m), and \( S \) are prognostic quantities in the model. Snow cover is assumed to be overlying vegetation and bare ground with the same fraction. The interception reservoir occupies an identical fraction of all snow-free tiles.

### Table 7.1  Vegetation Types and Parameter Values (See Text). H/L refer to the distinction between high and low vegetation.

<table>
<thead>
<tr>
<th>Index</th>
<th>Vegetation type</th>
<th>H/L</th>
<th>( \tau_{\text{sn}, \text{min}} ) (s m(^{-2}))</th>
<th>( \text{LAI} ) (m(^2)m(^{-2}))</th>
<th>( c_{\text{veg}} )</th>
<th>( g_0 ) (hPa (^{-1}))</th>
<th>( a_i )</th>
<th>( b_i )</th>
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<tbody>
<tr>
<td>1</td>
<td>Crops, mixed farming</td>
<td>L</td>
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<td>3</td>
<td>0.90</td>
<td>0</td>
<td>5.558</td>
<td>2.614</td>
</tr>
<tr>
<td>2</td>
<td>Short grass</td>
<td>L</td>
<td>110</td>
<td>2</td>
<td>0.85</td>
<td>0</td>
<td>10.739</td>
<td>2.608</td>
</tr>
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<td>3</td>
<td>Evergreen needleleaf trees</td>
<td>H</td>
<td>500</td>
<td>5</td>
<td>0.90</td>
<td>0.03</td>
<td>6.706</td>
<td>2.175</td>
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<td>Deciduous needleleaf trees</td>
<td>H</td>
<td>500</td>
<td>5</td>
<td>0.90</td>
<td>0.03</td>
<td>7.066</td>
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<td>Evergreen broadleaf trees</td>
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<td>0.03</td>
<td>5.990</td>
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<td>6</td>
<td>Deciduous broadleaf trees</td>
<td>H</td>
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<td>0.03</td>
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<td>1.303</td>
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<td>0</td>
<td>4.372</td>
<td>0.978</td>
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<td>-</td>
<td>-</td>
<td>-</td>
<td>-</td>
<td>-</td>
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<td>Bogs and marshes</td>
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<td>0</td>
<td>6.326</td>
<td>1.567</td>
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<td>0.03</td>
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<td>0.90</td>
<td>0.03</td>
<td>4.453</td>
<td>1.631</td>
</tr>
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<td>4</td>
<td>0.60</td>
<td>0</td>
<td>-</td>
<td>-</td>
</tr>
</tbody>
</table>
A small fraction \( f_{R} \) of net short-wave radiation that is transmitted directly to the top soil or snow layer. The remaining fraction of the short-wave radiation (1 - \( f_{R} \)) is absorbed by the skin layer.

Finally, the surface albedo, \( \alpha \), is similar for all land tiles within a grid box except for those covered with snow (see the snow scheme description below). The climate database provides the snow-free background albedo on a monthly basis. Long-wave emissivity, \( \varepsilon \), outside the window region is equal to 0.99 for all tiles; emissivity in the window region is tile dependent and varies between 0.93 and 0.98 (see Table 2.1 in Section 2.5.5 for more details). The remaining surface characteristics (roughness length for momentum, \( z_{int} \), and heat, \( z_{in} \)) are similar for all land tiles within a grid box and specified in the climate database (Chapter 10).
Table 7.2: Tile Specific Values.

<table>
<thead>
<tr>
<th>Index</th>
<th>Tile</th>
<th>$\Lambda_{sk}$ unstable (W m(^{-2})K(^{-1}))</th>
<th>$\Lambda_{sk}$ stable (W m(^{-2})K(^{-1}))</th>
<th>$fR_{i}$</th>
<th>Resistance scheme</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>Open water</td>
<td>$\infty$</td>
<td>$\infty$</td>
<td>0</td>
<td>Potential</td>
</tr>
<tr>
<td>2</td>
<td>Ice water</td>
<td>58</td>
<td>58</td>
<td>0</td>
<td>Potential</td>
</tr>
<tr>
<td>3</td>
<td>Interception reservoir</td>
<td>10</td>
<td>10</td>
<td>0.05</td>
<td>Resistance</td>
</tr>
<tr>
<td>4</td>
<td>Low vegetation</td>
<td>10</td>
<td>10</td>
<td>0.05</td>
<td>Resistance</td>
</tr>
<tr>
<td>5</td>
<td>Snow on low vegetation/bare ground</td>
<td>7</td>
<td>7</td>
<td>0</td>
<td>Potential</td>
</tr>
<tr>
<td>6</td>
<td>High vegetation</td>
<td>$\Lambda_{a,u} + 5$</td>
<td>$\Lambda_{a,s} + 5$</td>
<td>0.03</td>
<td>Resistance</td>
</tr>
<tr>
<td>7</td>
<td>High vegetation with snow beneath</td>
<td>$\Lambda_{a,u} + 5$</td>
<td>$\Lambda_{a,s} + 5$</td>
<td>0.03</td>
<td>Canopy and snow resistance</td>
</tr>
<tr>
<td>8</td>
<td>Bare ground</td>
<td>15</td>
<td>15</td>
<td>0</td>
<td>Resistance</td>
</tr>
</tbody>
</table>

The resistance scheme describes the way of coupling with the atmosphere: Potential denotes atmospheric resistance only; Resistance denotes aerodynamic resistance in series with a canopy or soil resistance; Canopy and snow resistance denotes a canopy resistance for the vegetation and an extra aerodynamic coupling to the snow surface (see Figs. 7.1–7.2 and Subsection 7.2.2). For tiles 6 and 7, $\Lambda_{a,u} = 15$ W m\(^{-2}\)K\(^{-1}\) and $\Lambda_{a,s} = 10$ W m\(^{-2}\)K\(^{-1}\) represent the aerodynamic coupling between the canopy and the soil in the unstable and stable cases, respectively, and the factor 5 represents the longwave radiative exchanges. Unstable/stable refers to the temperature gradient between the skin layer and the top soil or snow layer.

7.2.2 Surface heat and evaporation fluxes

A resistance parameterization is used to calculate the turbulent fluxes. Momentum exchange is parameterized with the same roughness length for all tiles, but with a different stability correction for each tile. The resistance scheme for water vapour and heat exchanges is different for different tiles (see Fig. 7.2). For ocean, sea ice and snow on low vegetation, the turbulent fluxes of heat and water vapour are given by

\[
H_i = \rho_a c_p |U_L| C_{H,i} (T_L + g z_L - T_{sk,i})
\]

(7.5)

\[
E_i = \rho_a |U_L| C_{H,i} [q_L - q_{sat}(T_{sk,i})]
\]

(7.6)

with $\rho_a$ the air density, $c_p$ the heat capacity of moist air, $g$ the acceleration of gravity, $|U_L|$, $T_L$, $q_L$, $z_L$ the wind speed, temperature, humidity and height of the lowest atmospheric model level, and $C_{H,i}$ the turbulent exchange coefficient, that varies from tile to tile because of different atmospheric stabilities. See Chapter 3 for a description of the exchange coefficients where different roughness lengths for heat and momentum are assumed and a Monin–Obukhov formulation is adopted for the stability dependence.
Figure 7.2  Resistance scheme for three categories of coupling. Potential refers to ocean, sea ice and and snow on low vegetation; (Canopy) resistance to dry low and dry high vegetation, bare soil, and interception reservoir when potential evaporation exceeds the maximum reservoir content; Resistance to snow to snow under high vegetation.

For high and low vegetation, an additional canopy resistance $r_c$ is added:

$$E_i = \frac{P_s}{r_a + r_c} [q_L - q_{sat}(T_{sk,i})]$$  \hspace{1cm} (7.7)

with $r_a = (|U_L| C_{l,i})^{-1}$ and $i$ indicating the high or low vegetation tiles. $r_c$ is a function of downward shortwave radiation $R_c$, leaf area index $LAI$, average unfrozen root soil water $\bar{\theta}$, atmospheric water vapour deficit $D_a$ and a minimum stomatal resistance $r_{s,\text{min}}$, following Jarvis (1976):

$$r_c = \frac{r_{s,\text{min}}}{LAI} f_1(R_c) f_2(\bar{\theta}) f_3(D_a)$$  \hspace{1cm} (7.8)

$f_1$ is a hyperbolic function of downward short-wave radiation only:

$$\frac{1}{f_1(R_c)} = \min \left[ \frac{bR_c + c}{a(bR_c + 1)} \right]$$  \hspace{1cm} (7.9)

where $a = 0.81$, $b = 0.004$ W m$^{-2}$ and $c = 0.05$.

Function $f_2$ is defined as

$$\frac{1}{f_2(\bar{\theta})} = \begin{cases} 0 & \bar{\theta} < \theta_{\text{pwp}} \\ \bar{\theta} - \theta_{\text{pwp}} & \theta_{\text{pwp}} \leq \bar{\theta} \leq \theta_{\text{cap}} \\ \theta_{\text{cap}} - \theta_{\text{pwp}} & \theta_{\text{pwp}} \leq \bar{\theta} \leq \theta_{\text{cap}} \\ 0 & \bar{\theta} > \theta_{\text{cap}} \end{cases}$$  \hspace{1cm} (7.10)

where the soil moisture at permanent wilting point and at field capacity, $\theta_{\text{pwp}}$ and $\theta_{\text{cap}}$, respectively, are defined...
in Table 7.5, $\bar{\theta}$ is a weighted average of the unfrozen soil water

$$\bar{\theta} = \sum_{k=1}^{\infty} R_k \max[\theta_{\text{liq},k} \theta_{\text{wp},k}]$$

(7.11)

where $R_k$ is the fraction of roots in layer $k$ and the fraction of unfrozen soil water, $f_{\text{liq},k} = 1 - f_{\text{liq}}(T_k)$, is a parameterized function of the soil temperature of layer $k$, $T_k$, as specified in Section 7.5.2. Table 7.1 lists the coefficients $a_i$ and $b_i$ which are used to calculate the root fraction $R_k$ according to Zeng et al. (1998):

$$R_k = 0.5 \{ \exp(-a_k z_k^{-1/2}) + \exp(-b_k z_k^{-1/2}) - \exp(-a_k z_k^{+1/2}) - \exp(-b_k z_k^{+1/2}) \}$$

(7.12)

where $z_{k+1/2}$ is the depth of the bottom of layer $k$ (in m; $z_{1/2} = 0$ m). Contributions from levels exceeding the column depth are added to the deepest soil layer in order to ensure that $\sum R_k = 1$. Table 7.3 lists the distribution of the roots over the four soil layers.

<table>
<thead>
<tr>
<th>Vegetation index</th>
<th>1</th>
<th>2</th>
<th>3</th>
<th>4</th>
<th>5</th>
<th>6</th>
<th>7</th>
<th>8</th>
<th>9</th>
<th>10</th>
<th>11</th>
<th>12</th>
<th>13</th>
<th>14</th>
<th>15</th>
</tr>
</thead>
<tbody>
<tr>
<td>Layer 1</td>
<td>24</td>
<td>35</td>
<td>26</td>
<td>26</td>
<td>24</td>
<td>25</td>
<td>27</td>
<td>100</td>
<td>47</td>
<td>24</td>
<td>17</td>
<td>25</td>
<td>23</td>
<td>23</td>
<td>19</td>
</tr>
<tr>
<td>Layer 2</td>
<td>41</td>
<td>38</td>
<td>39</td>
<td>38</td>
<td>38</td>
<td>34</td>
<td>27</td>
<td>0</td>
<td>45</td>
<td>41</td>
<td>31</td>
<td>34</td>
<td>36</td>
<td>36</td>
<td>35</td>
</tr>
<tr>
<td>Layer 3</td>
<td>31</td>
<td>23</td>
<td>29</td>
<td>29</td>
<td>31</td>
<td>27</td>
<td>27</td>
<td>0</td>
<td>8</td>
<td>31</td>
<td>33</td>
<td>27</td>
<td>30</td>
<td>30</td>
<td>36</td>
</tr>
<tr>
<td>Layer 4</td>
<td>4</td>
<td>4</td>
<td>6</td>
<td>7</td>
<td>7</td>
<td>14</td>
<td>9</td>
<td>0</td>
<td>0</td>
<td>4</td>
<td>19</td>
<td>11</td>
<td>11</td>
<td>10</td>
<td>10</td>
</tr>
</tbody>
</table>

TABLE 7.3 ROOT DISTRIBUTION PER VEGETATION TYPE (IN %) OVER THE FOUR LAYERS. VEGETATION INDEXES REFER TO TABLE 7.1.

A dependence on atmospheric humidity deficit ($D = e_{\text{sat}}(T) - e_L$, with $e$ the vapour pressure) is included according to

$$\frac{1}{f_3(D_a)} = \exp(-g_D D_a)$$

(7.13)

where $g_D$ depends on the vegetation type (Table 7.1), and is non-zero for high vegetation only.

Evaporation from the interception reservoir is given by Eq. (7.6) only when the amount of water in the interception reservoir, $W_I$, is sufficient to sustain potential evaporation during the entire time step $\Delta t$. If $W_I$ is limited, an additional resistance $r_I$, analogue to $r_e$ in Eq. (7.7), is introduced. $r_I$ is calculated from the potential evaporation of the previous time step. Note that this type of flux-limiter is a time-step dependent feature of the model numerics.

Bare-soil evaporation uses a resistance approach, an analogue to the canopy transpiration formulation (Eq. (7.7)). The soil evaporation resistance, $r_{\text{soil}}$, is

$$r_{\text{soil}} = r_{\text{soil, min}} f_2(f_{\text{liq}} \theta_1)$$

(7.14)

with $f_2$ given by Eq. (7.10), and $r_{\text{soil, min}} = 50$ s m$^{-1}$. By this parameterization, evaporation from bare ground is treated similar to a single leaved canopy with a minimum resistance $r_{\text{soil, min}}$ extracting water from the upper soil layer only, and not experiencing any additional stress due to limited radiation or dry air. Eq. (7.14) shuts off evap-
oration when the top soil moisture reaches permanent wilting point. When compared to observations over semi-arid areas, an alternative relative humidity formulation (Mahrour and Noilhan 1991; Viterbo and Belskis 1995), that does not have a similar limitation, gave excessive evaporation (van den Hurk et al. 2000).

A special treatment is included in the calculation of evaporation over high vegetation with snow underneath (see Fig. 7.2). Evaporation takes place from both the canopy component in the tile \( E_{\text{veg},7} \) and from the snow lying under the vegetation. The canopy evaporation uses a canopy resistance and saturation specific humidity at the canopy skin temperature, while the snow evaporation \( E_{\text{sn},7} \) is parameterized with an additional constant aerodynamic resistance \( r_{a,\text{sn}} \) and saturation specific humidity at snow temperature \( T_{\text{sn}} \). The evaporation from tile 7 is the combination of the canopy transpiration and the snow evaporation:

\[
E_7 = \rho_a \frac{q_L - q_*}{r_a} = \frac{q_* - q_{\text{sat}}(T_{\text{sn}})}{r_{a,\text{sn}}} + \frac{q_* - q_{\text{sat}}(T_{\text{sk}})}{r_c}
\]

(7.15)

where \( q_* \) is the humidity at the connection point of the three resistances (Fig. 7.2). After elimination of \( q_* \), \( E_7 \) can be rewritten as:

\[
E_7 = \rho_a \frac{q_L - q_{\text{sat}}(T_{\text{sk}})}{r_a + \frac{r_a}{r_{a,\text{sn}}}} + \rho_a \frac{q_* - q_{\text{sat}}(T_{\text{sn}})}{r_a + r_c + \frac{r_a}{r_{a,\text{sn}}}} \frac{r_a}{r_{a,\text{sn}} + r_{a,\text{sn}}r_c}
\]

(7.16)

The first term in the equation above is interpreted as \( E_{\text{veg},7} \) and is treated in the standard way (i.e., implicit in the tile skin temperature). The second term is interpreted as evaporation from snow \( E_{\text{sn},7} \) and is handled explicitly. The values of \( r_{a,\text{sn}} \) depend on the stability of the subcanopy layer and are functions of \( \Lambda_{a,u} \) and \( \Lambda_{a,s} \) (see Table 7.2); \( r_{a,\text{sn}} = 67 \text{ s m}^{-1} \) and \( r_{a,\text{sn}} = 220 \text{ s m}^{-1} \) for an unstable and stable subcanopy layer, respectively. In spring, the latent heat flux of that tile, \( L_v E_{\text{veg},7} + L_e E_{\text{sn},7} \) will be dominated by snow evaporation since the frozen soil under the snow deck will lead to very large values of \( r_c \).

The grid box total sensible and latent heat fluxes are expressed as an area weighted average:

\[
H = \sum_{i=1}^{8} C_i H_i
\]

(7.17)

\[
E = \sum_{i=1}^{8} C_i E_i
\]

(7.18)

with \( H_i \) given by Eq. (7.5), and \( E_i \) by Eq. (7.6) for ocean, sea-ice and snow on low vegetation, Eq. (7.7) for dry high and low vegetation, the interception reservoir (with \( r_c \) replaced by \( r_l \)) and for bare soil (with \( r_c \) replaced by \( r_{\text{soil}} \)) and Eq. (7.16) for high vegetation with underlying snow.

### 7.3 The Surface Energy Balance and Coupling to the Soil

A skin temperature \( T_{sk} \) forms the interface between the soil and the atmosphere. As detailed in Section 3.5, it is calculated for each grid box tiles separately, by scanning the surface energy balance solver over the 8 tiles, assuming a complete coverage of the specific tile. For a single tile, this procedure is very similar to the derivation of the Penman–Monteith equation in which the skin temperature is eliminated from the surface energy balance equation.
The numerical approach used in TESSELS has the advantage that the feedback of skin temperature on net radiation and ground heat flux is included (see Section 3.5). The input radiation and reference atmospheric temperature ($T_L$), specific humidity ($q_L$), and wind speed ($U_L$) are identical for each tile. The surface fluxes “seen” by the atmosphere are calculated as an area-weighted average over the tiles (see Eqs. (7.17) and (7.18)). For the high vegetation with snow underneath, the skin temperature is that of the high vegetation; the temperature of the underlying snow is calculated separately.

The energy balance equation solved for each tile takes into account partial absorption of net short-wave radiation, $1-f_{RS,i}$, in the skin layer (see Table 7.2). The remaining energy is directly passed to the soil or snow:

$$(1-f_{RS,i})(1-\alpha_i)R_s + \varepsilon(R_T-\sigma T_{sk,i}^4) + H_i + L_{vs,i}E_i = \Lambda_{sk,i}(T_{sk,i} - T_1)$$

(7.19)

where $i$ denotes the tile index, $R_s$ and $R_T$ are downward short-wave radiation and long-wave radiation, respectively, $\sigma$ is the Stefan–Boltzmann constant, $T_1$ the temperature of the upper soil or snow layer, $H_i$ the sensible heat flux, and $L_{vs,i}E_i$ the latent heat flux from the skin layer, and $\Lambda_{sk,i}$, the skin conductivity for tile $i$. Latent heat of evaporation, $L_v$, is used for all evaporation terms except snow evaporation, while $L_s$, the latent heat of sublimation, is used for evaporation of snow (i.e., tile 5 and the contribution $E_{sn,7}$ from tile 7, defined by Eq. (7.16)).

The tiled surface is thermally coupled to the snow deck, when present, and to a single soil profile. The net flux into the soil is a weighted average of the flux from each tile.

The solution of Eq. (7.19) is performed inside the code for turbulent exchanges in the atmosphere (Chapter 3). The atmospheric vertical diffusion equations yield a tridiagonal system of equations, with the coupling to the skin temperature given by the matrix row corresponding to the lowest model level. The first step for the solution of the system of equations, an LU decomposition, is followed by the solution of Eq. (7.19) before back-substitution. Details of the computations can be found in Chapter 3.

### 7.4 Snow

The snow scheme represents an additional “layer” on top of the upper soil layer, with an independent, prognostic, thermal and mass contents. The snow pack is represented by a single snow temperature, $T_{sn}$, and the snow mass per unit area (snow mass for short) $S$. The net energy flux at the top of the snow pack, $G_{sn}^T$, is the residual of the skin energy balance from the snow covered tiles and the snow evaporation from the tile with high vegetation over snow (Eq. (7.15)). The basal heat flux, $G_{sn}^B$, is given by equation a resistance formulation modified in case of melting. The absorbed energy is used to change the snow temperature or melt the snow, when $T_{sn}$ exceeds the melting point.

The heat capacity of the snow deck is a function of its depth and the snow density, which is a prognostic quantity depending on snow age following (Douville et al. 1995). The snow thermal conductivity changes with changing snow density. The snow albedo changes exponentially with snow age. For snow on low vegetation it ranges between 0.50 for old snow and 0.85 for fresh snow (to which it is reset whenever the snow fall exceeds 1 mm hr$^{-1}$). The albedo for high vegetation with snow underneath is fixed at 0.15.

#### 7.4.1 Snow mass and energy budget

The snow mass budget reads as:

$$\rho_s \frac{\partial S}{\partial t} = F + c_{sn}(E_{sn} - M_{sn})$$

(7.20)
Chapter 7 ‘Surface parametrization’

where \( F \) is snowfall (units kg m\(^{-2}\)), \( S \) is snow mass (sometimes referred as snow water equivalent) grid-averaged (units 103 kg m\(^{-2}\)), \( \rho_w \) is the water density (units kg m\(^{-3}\)), \( E_{sn} \) and \( M_{sn} \) are snow evaporation and melting, respectively (units kg m\(^{-3}\)), and \( c_{sn} \) is the snow fraction (see Eq. (7.2)), i.e. the sum of tiles 5 and 7 (see Eq. (7.4)). In Eq. (7.20) and in the remaining of this section, all surface fluxes are per unit area and apply only to the snow area (i.e. tile 5 and 7). The snow equivalent water \( S \) applies to the entire grid square and therefore occurs in the equation divided by the total snow fraction. The snow flux from the atmospheric model, \( F \), is again for the entire grid square. As a general rule, all quantities with subscript \( sn \) will refer to the snow area. In Eq. (7.20), the snow evaporation is defined as

\[
e_{sn}E_{sn} = e_{5}E_{5} + c_{7}E_{sn,7} \tag{7.21}
\]

Snow mass and snow depth are related by

\[
D_{sn} = \frac{\rho_w S}{\rho_{sn} c_{sn}} \tag{7.22}
\]

where \( D_{sn} \) is snow depth for the snow-covered area (units m; \( D_{sn} \) is NOT a grid-averaged quantity) and \( \rho_{sn} \) is the snow density (units kg m\(^{-3}\)).

The snow energy budget reads as

\[
(pC)_{sn}D_{sn} \frac{dT_{sn}}{dt} = (pC)_{sn} \frac{\rho_w S}{\rho_{sn} c_{sn}} \frac{dT_{sn}}{dt} = (pC)_{sn} \frac{\rho_w S}{\rho_{sn} c_{sn}} \frac{dT_{sn}}{dt} = R_{sn}^N + L_{sn}E_{sn} + H_{sn} - G_{sn}^B - Q_{sn} \tag{7.23}
\]

where \((pC)_{sn}\) and \((pC)_{i}\) are the ice and snow volumetric heat capacities, respectively (units J m\(^{-3}\)K\(^{-1}\)), \( \rho_i \) is the ice density (units kg m\(^{-3}\)), \( R_{sn}^N \) is the net radiation absorbed by the snow pack (units W m\(^{-2}\)), \( L_i \) is the latent heat of sublimation (units J kg\(^{-1}\)), \( H_{sn} \), \( G_{sn}^B \), and \( Q_{sn} \) represent, respectively, the snow sensible heat flux, basal heat flux (at the bottom of the snow pack), and energy exchanges due to melting (units W m\(^{-2}\)). Eq. (7.23) neglects the thermal energy brought by precipitation. The snow is composed of an ice fraction, a liquid water fraction and an air fraction, \( v_i \), \( v_w \) and \( v_a \), respectively, where typically \( 0.3 < v_a < 0.9 \) and the liquid water fraction is significantly different from zero in melting conditions. The following approximations are made in Eq. (7.23)

\[
\frac{(pC)_{sn}}{\rho_{sn}} = \frac{(pv_i) + v_w(pv) + v_a(pC)}{v_i + v_w + v_a} = \frac{(pv_i) + v_w(pv) + v_a(pC)}{v_i + v_w + v_a} = \frac{(pv_i) + v_w(pv) + v_a(pC)}{v_i + v_w + v_a} \tag{7.24}
\]

The melting term couples the mass and energy equation

\[
Q_{sn} = L_{m}M_{sn} = -L_{m} \left. \frac{\rho_w S}{c_{sn}} \right|_{\text{m}} \tag{7.25}
\]

where \( L_m \) is the latent heat of fusion (units J kg\(^{-1}\)) and the subscript \( m \) represents melting.

### 7.4.2 Prognostic snow density and albedo

Following Douville et al. (1995) snow density is assumed to be constant with depth and to evolve exponentially towards a maximum density (Verseghy, 1991). First a weighted average is taken between the current density and
the minimum density for fresh snow

\[ \rho_{sn}^* = \frac{S \rho_{sn}^f + (\Delta t F/\rho_w) \rho_{min}}{S + (\Delta t F/\rho_w)} \]  

(7.26)

The exponential relaxation reads

\[ \rho_{sn}^{t+1} = (\rho_{sn}^* - \rho_{max}) \exp(-\tau_1 \Delta t/\tau_t) + \rho_{max} \]  

(7.27)

where timescales \( \tau_1 = 86400 \) s, and \( \tau_t = 0.24 \) corresponding to an e-folding time of about 4 days, with minimum density \( \rho_{min} = 100 \) kg m\(^{-3}\) and maximum density \( \rho_{max} = 300 \) kg m\(^{-3}\) (see Table 7.4).

<table>
<thead>
<tr>
<th>Symbol</th>
<th>Parameter</th>
<th>Value</th>
</tr>
</thead>
<tbody>
<tr>
<td>( D_{sn}^{max} )</td>
<td>Maximum snow thermal depth</td>
<td>0.07 m</td>
</tr>
<tr>
<td>( S_{cr} )</td>
<td>Threshold value for grid box coverage of snow</td>
<td>0.015 m</td>
</tr>
<tr>
<td>( \alpha_{min} )</td>
<td>Minimum albedo of exposed snow</td>
<td>0.50</td>
</tr>
<tr>
<td>( \alpha_{max} )</td>
<td>Maximum albedo of exposed snow</td>
<td>0.85</td>
</tr>
<tr>
<td>( \alpha_{sn,f} )</td>
<td>Albedo of shaded snow</td>
<td>0.2</td>
</tr>
<tr>
<td>( \lambda_i )</td>
<td>Ice heat conductivity</td>
<td>2.2 W m(^{-1})K(^{-1})</td>
</tr>
<tr>
<td>( \rho_{min} )</td>
<td>Minimum snow density</td>
<td>300 k gm(^{-3})</td>
</tr>
<tr>
<td>( \rho_{max} )</td>
<td>Maximum snow density</td>
<td>100 k gm(^{-3})</td>
</tr>
<tr>
<td>( \rho_i )</td>
<td>Ice density</td>
<td>920 kg m(^{-3})</td>
</tr>
<tr>
<td>( (\rho C)_i )</td>
<td>Ice volumetric heat capacity</td>
<td>2.05 ( 10^6 ) J m(^{-3}) K(^{-1})</td>
</tr>
<tr>
<td>( \tau_a )</td>
<td>Linear coefficient for decrease of albedo of non-melting snow</td>
<td>0.008</td>
</tr>
<tr>
<td>( \tau_1 )</td>
<td>Coefficient for exponential decrease of snow density and melting snow albedo</td>
<td>0.24</td>
</tr>
<tr>
<td>( \tau_j )</td>
<td>Length of day</td>
<td>86400 s</td>
</tr>
</tbody>
</table>

Snow albedo in exposed areas evolves according to the formulation of Baker et al. (1990), Verseghy (1991) and Douville et al. (1995). For non melting-conditions:

\[ \alpha_{sn}^{t+1} = \alpha_{sn}^f - \tau_1 \Delta t/\tau_1 \]  

(7.28)

where \( \tau_a = 0.008 \), which will decrease the albedo by 0.1 in 12.5 days. For melting conditions \( M_{sn} > 0 \):

\[ \alpha_{sn}^{t+1} = (\alpha_{sn}^f - \alpha_{min}) \exp(-\tau_1 \Delta t/\tau_1) + \alpha_{min} \]  

(7.29)

where \( \alpha_{min} = 0.5 \) and \( \alpha_{max} = 0.85 \). If snowfall \( F > 1 \) kg m\(^{-3}\)hr\(^{-1}\), the snow albedo is reset to the maximum value, \( \alpha_{sn}^{t+1} = \alpha_{max} \).

The above formulae are inadequate to describe the evolution of the surface albedo of snow cover with high vegetation. Observations suggest a dependence on forest type but, by and large, the albedo changes from a value around
0.3 just after a heavy snowfall to a value around 0.2 after a few days (see Betts and Ball (1997) and the discussion in Viterbo and Betts (1999)). This change reflects the disappearance of intercepted snow, due to melt (for sufficiently warm temperatures) or wind drift (for cold temperatures). Ways of describing those two mechanisms would involve either a separate albedo variable for the snow in the presence of high vegetation, or the introduction of an interception reservoir for snow. In the absence of any of the two, we define $\alpha_{sn,f} = 0.2$ for the snow in the presence of high vegetation. This value was chosen to match the overall forest albedo in the presence of snow from the results of Viterbo and Betts (1999).

### 7.4.3 Additional details

#### 7.4.3 (a) Limiting of snow depth in the snow energy equation

Initial experimentation with the snow model revealed that the time evolution of snow temperature was very slow over Antarctica. The reason is rather obvious; the snow depth over Antarctica is set to a climatological value of 10 m which can respond only very slowly to the atmospheric forcing due to its large thermal inertia. In previous model versions, the properties of layer 1 were replaced by snow properties when snow was present, which kept the timescale short. A physical solution would have been to introduce a multilayer snow model, with e.g. four layers to represent timescales from one day to a full annual cycle. As a shortcut, a limit is put on the depth of the snow layer in the thermal budget, $D_{sn}^{max} = 0.07$ m. The energy equation reads:

$$
(pC)_{sn}D_{sn}^{*} \frac{\partial T_{sn}}{\partial t} = R_{sn}^{N} + L_{sn}E_{sn} + H_{sn} - C_{sn}^{B} - Q_{sn}
$$

$$D_{sn}^{*} = \min(D_{sn}, D_{sn}^{max})$$

(7.30)

#### 7.4.3 (b) Basal heat flux and thermal coefficients

The heat flux at the bottom of the snow pack is written as a finite difference in the following way:

$$G_{sn}^{B} = \frac{T_{sn} - T_{1}}{r_{sn}}$$

(7.31)

where $r_{sn}$ is the resistance between the middle of the snow pack and the middle of soil layer 1, with two components: the resistance of the lower part of the snow pack and the resistance of the top half of soil layer 1:

$$r_{sn} = 0.5 \frac{D_{sn}^{*}}{\lambda_{sn}} + \frac{1}{\Lambda_{sk,8}}$$

(7.32)

where the second term is the skin layer conductivity for bare soil (tile 8), which can be seen as an approximation of $0.5(D_{1}/\lambda_{r})$. The snow thermal conductivity, is related to the ice thermal conductivity according to Douville et al. (1995):

$$\lambda_{sn} = \lambda\left(\frac{\rho_{sn}}{\rho_{i}}\right)^{1.88}$$

(7.33)

Table 7.4 contains the numerical values of the ice density and ice heat conductivity.

#### 7.4.3 (c) Numerical solution for non-melting situations

The net heat flux that goes into the top of the snow deck is an output of the vertical diffusion scheme
\[ H_{sn}^N = R_{sn}^N + L_{sub}E_{sn} + H_{sn} \quad (7.34) \]

In the absence of melting, the solution of Eq. (7.30) is done implicitly. The preliminary snow temperature, prior to the checking for melting conditions, \( T_{sn}^* \), is given by

\[ A_1 \frac{T_{sn}^* - T_{sn}^t}{\Delta t} = H_{sn}^N - \frac{T_{sn}^* - T_1}{r_{sn}} \quad (7.35) \]

\[ A_1 = \min \left[ (\rho C_i) \frac{\rho_w}{\rho_l} \frac{S}{c_{sn}}, A_1^{\text{max}} \right] \]

\[ A_1^{\text{max}} = (\rho C_i) \frac{\rho_w}{\rho_l} D_{sn}^{\text{max}} \quad (7.36) \]

where superscript \( t \) refers to the current time step and superscript * to the preliminary value at the next time step.

The solution for \( T_{sn}^* \) is obtained from

\[ T_{sn}^* \left( 1 + \frac{\Delta t}{r_{sn} A_1} \right) = T_{sn}^t + \frac{\Delta t}{A_1} \left( H_{sn}^N + \frac{T_1}{r_{sn}} \right) \quad (7.37) \]

The basal snow heat flux to be used as input for the thermal budget of the soil (in the snow covered fraction only) is

\[ G_{sn}^B = \frac{T_{sn}^* - T_1}{r_{sn}} \quad (7.38) \]

Finally, a preliminary new value for the snow mass, \( S^* \), is computed from snow fall and snow evaporation

\[ \rho_w \frac{S^* - S^t}{\Delta t} = F + c_{sn} E_{sn} \quad (7.39) \]

### 7.4.4 Treatment of melting

**7.4.4 (a) No melting occurs.** If \( T_{sn}^* < T_0 \), no melting occurs and the preliminary values \( T_{sn}^* \) and \( S^* \) become the \( t+1 \) values, while the basal heat flux is given by Eq. (7.38).

**7.4.4 (b) Melting conditions.** If \( T_{sn}^* > T_0 \), snow melting occurs and the time step is divided in two fractions, \( \Delta t = \Delta_1 t + \Delta_2 t \), where the first fraction, \( \Delta_1 t \) brings the temperature to \( T_0 \) with no melting:

\[ \Delta_1 t = \frac{A_1 (T_0 - T_{sn}^t)}{H_{sn}^N - (T_0 - T_1)/r_{sn}} \quad (7.40) \]

while, during the second fraction, \( \Delta_2 t \), melting occurs with no resultant warming of the snow:
\[ T_{n+1} = T_0 \]
\[ Q_{sn} = H_{sn}^N - G_{sn}^B \]
\[ \rho_s \Delta_2 t \frac{S_{n+1} - S^n}{H_{sn}^N - G_{sn}^B} = -c_s M_{sn} = -\frac{Q_{sn}}{L_f} = -c_s \frac{H_{sn}^N - G_{sn}^B}{L_f} \]

(7.41)

**If not all the snow melts**, i.e., if \( S_{n+1} > 0 \), the following heat flux is passed to the soil

\[ G_{sn}^B = \frac{T_0 - T_1}{T_{sn}^B} \]

(7.42)

**When all the snow melts**, i.e., if \( S_{n+1} < 0 \), the melting time step is redefined as:

\[ S_{n+1} = 0 \]
\[ \Delta_2 t = \rho_s L_f \frac{S^n}{c_s (H_{sn}^N - G_{sn}^B)} \]
\[ \Delta_3 t = 1 - (\Delta_1 t + \Delta_2 t) \]

and the basal heat flux is redefined as

\[ G_{sn}^B = \frac{\Delta_1 t + \Delta_2 t}{t_{sn}} \frac{T_0 - T_1}{T_{sn}^B} + \frac{\Delta_3 t}{\Delta t} \frac{H_{sn}^N}{L_f} \]

(7.44)

### 7.5 Soil Heat Transfer

In the absence of internal phase changes, the soil heat transfer is assumed to obey the following Fourier law of diffusion

\[ (\rho C)_{soil} \frac{\partial T}{\partial t} = \frac{\partial}{\partial z} \left[ \lambda_T \frac{\partial T}{\partial z} \right] \]

(7.45)

where \((\rho C)_{soil}\) is the volumetric soil heat capacity \( (\text{J m}^{-3}\text{K}^{-1}) \), \( T \) is the soil temperature (units K), \( z \) is the vertical coordinate—the distance from the surface, positive downwards—(units m), and \( \lambda_T \) is the thermal conductivity \( (\text{W m}^{-1}\text{K}^{-1}) \). The above equation assumes that heat fluxes are predominantly in the vertical direction, that the effects of phase changes in the soil and the heat transfer associated with the vertical movement of water in the soil can be neglected (de Vries 1975), and that the effects of hysteresis can be neglected (Milly 1982).

The boundary condition at the bottom, no heat flux of energy, is an acceptable approximation provided that the total soil depth is large enough for the time-scales represented by the model or, in other words, the bottom of the soil is specified at a depth where the amplitude of the soil heat wave is a negligible fraction of its surface amplitude (see de Vries (1975) and next section).
7.5.1 Discretization and choice of parameters

**Table 7.5** Parameters in the land-surface scheme. See Table 7.4 for snow-related parameters.

<table>
<thead>
<tr>
<th>Symbol</th>
<th>Parameter</th>
<th>Value</th>
</tr>
</thead>
<tbody>
<tr>
<td>( b )</td>
<td>Clapp and Hornberger soil parameter</td>
<td>6.04</td>
</tr>
<tr>
<td>( b_I )</td>
<td>Interception efficiency</td>
<td>0.25</td>
</tr>
<tr>
<td>( D_1 )</td>
<td>Depth of soil layer 1</td>
<td>0.07 m</td>
</tr>
<tr>
<td>( D_2 )</td>
<td>Depth of soil layer 2</td>
<td>0.21 m</td>
</tr>
<tr>
<td>( D_3 )</td>
<td>Depth of soil layer 3</td>
<td>0.72 m</td>
</tr>
<tr>
<td>( D_4 )</td>
<td>Depth of soil layer 4</td>
<td>1.89 m</td>
</tr>
<tr>
<td>( F_{cv} )</td>
<td>Fraction of gridbox covered by convective rainfall</td>
<td>0.5</td>
</tr>
<tr>
<td>( k )</td>
<td>Heterogeneity factor for convective precipitation</td>
<td>0.5</td>
</tr>
<tr>
<td>( T_{11} )</td>
<td>Highest temperature for existence of ice water</td>
<td>( T_0 + 1 )</td>
</tr>
<tr>
<td>( T_{12} )</td>
<td>Lowest temperature for existence of liquid water</td>
<td>( T_0 - 3 )</td>
</tr>
<tr>
<td>( W_{\text{max}} )</td>
<td>Maximum water amount on single leaf</td>
<td>0.0002 m</td>
</tr>
<tr>
<td>( \gamma_{\text{sat}} )</td>
<td>Hydraulic conductivity at saturation</td>
<td>( 4.57 \times 10^{-4} ) m s(^{-1} )</td>
</tr>
<tr>
<td>( \lambda_{\text{dry}} )</td>
<td>Heat conductivity of dry soil</td>
<td>0.190 W m(^{-1})K(^{-1})</td>
</tr>
<tr>
<td>( \lambda_{\text{sm}} )</td>
<td>Heat conductivity of soil matrix</td>
<td>3.44 W m(^{-1})K(^{-1})</td>
</tr>
<tr>
<td>( \lambda_{\text{w}} )</td>
<td>Heat conductivity of liquid water</td>
<td>0.57 W m(^{-1})K(^{-1})</td>
</tr>
<tr>
<td>((\rho C)_{\text{soil}})</td>
<td>Volumetric soil heat capacity</td>
<td>( 2.19 \times 10^6 ) J m(^{-3})K(^{-1})</td>
</tr>
<tr>
<td>( \theta_{\text{sat}} )</td>
<td>Soil moisture at saturation</td>
<td>0.472 m(^3)m(^{-3})</td>
</tr>
<tr>
<td>( \theta_{\text{cap}} )</td>
<td>Soil moisture at field capacity</td>
<td>0.323 m(^3)m(^{-3})</td>
</tr>
<tr>
<td>( \theta_{\text{pwp}} )</td>
<td>Soil moisture at permanent wilting point</td>
<td>0.171 m(^3)m(^{-3})</td>
</tr>
<tr>
<td>( \psi_{\text{sat}} )</td>
<td>Matric potential at saturation</td>
<td>(-0.338) m</td>
</tr>
</tbody>
</table>

For the solution of Eq. (7.45) the soil is discretized in four layers, of depths \( D_k \), \( k = 1, 2, 3, 4 \), the temperatures are defined at full layers \( T_k \), and the heat fluxes, at half layers (\( G_{k+1/2} \)) is the heat flux, positive downwards, units W m\(^{-2}\), at the interface between layer \( k \) and \( k + 1 \). An energy-conserving implicit algorithm is used, leading to a tridiagonal system of equations with solution detailed in Section 7.8.

The boundary condition at the bottom is:

\[
G_{4+1/2} = 0
\]  

(7.46)

At the top, the boundary condition is the soil heat flux at the surface, computed as a weighted average over the tiles. For the snow free tiles, the flux into the soil consists of two parts. Apart from the diffusion of heat governed by \( \Lambda_{k,i}(T_{sk,i} - T_i) \) (see Eq. (7.19)), the net shortwave radiation not absorbed by the skin layer \( f_{s,i} \) provides energy to the soil. \( \text{Table 7.2} \) lists the values of \( \Lambda_{k,i} \) and \( f_{s,i} \) for each of the tiles. For the snow tiles, the heat flux into the soil is the snow basal flux, calculated using a resistance formulation and modified in the case of partial melting (see Eqs. (7.31), (7.38), (7.42), and (7.44)).

The net heat flux into the soil is given by:
where the summation scans all snow free tiles.

The volumetric soil heat capacity is assumed constant, with value \(2.19 \times 10^6\) J m\(^{-3}\) K\(^{-1}\) (see Table 7.5 for a list of constants used by the model). The heat conductivity, \(\lambda\), depends on the soil-water content following Peters-Lidard et al. (1998) (see also Farouki 1986; Johansen 1975) and is given by a combination of dry \(\lambda_{\text{dry}}\) and saturated \(\lambda_{\text{sat}}\) values, weighted by a factor known as the Kersten number, \(K_e\):

\[
\lambda = K_e (\lambda_{\text{sat}} - \lambda_{\text{dry}}) + \lambda_{\text{dry}},
\]

(7.48)

where \(\lambda_{\text{dry}} = 0.190\) W m\(^{-1}\) K\(^{-1}\) and

\[
\lambda_{\text{sat}} = \lambda_{\text{sm}}^{1 - \theta_{\text{sat}}} \lambda_w^{\theta_{\text{sat}}},
\]

(7.49)

where the heat conductivity of the soil matrix, \(\lambda_{\text{sm}} = 3.44\) W m\(^{-1}\) K\(^{-1}\) and the thermal conductivity of water is \(\lambda_w = 0.57\) W m\(^{-1}\) K\(^{-1}\). Eq. (7.49) represents a simplification of Peters-Lidard formulation, neglecting the changes in conductivity due to ice water and assuming the quartz content typical of a loamy soil. Finally, the Kersten number for fine soils was selected in Peters-Lidard et al. (1998):

\[
K_e = \log_{10} \left[ \max \left( 0.1; \frac{\theta_{\text{sat}}}{\theta_{\text{sat}}} \right) \right] + 1
\]

(7.50)

The depths of the soil layers are chosen in an approximate geometric relation (see Table 7.5), as suggested in Dardaroff (1978). Warrilow et al. (1986) have shown that four layers are enough for representing correctly all timescales from one day to one year. Using the numerical values of the heat capacity and soil depths defined in Table 7.5, the amplitude and phase response of the numerical solution of Eq. (7.45) were analysed by Viterbo and Beljaars (1995) for typical values of soil moisture in Eq. (7.48), and for harmonic forcings at the surface with periods ranging from half a day to two years. The analysis points to an error in the numerical solution of less than 20% in amplitude and 5% in phase for forcing periods between one day and one year.

### 7.5.2 Soil-water phase changes

At high and mid latitudes the phase changes of water in the soil have an important effect on the water and energy transfer in the soil. A proper consideration of the solid phase of soil water requires modifications including, in order of importance:

- **(a)** The thermal effects related to the latent heat of fusion/freezing (e.g. Rouse 1984);
- **(b)** Changes in the soil thermal conductivity due to the presence of ice (e.g. Penner 1970, not included in TESSEL as mentioned in the previous section);
- **(c)** Suppression of transpiration in the presence of frozen ground (e.g. Betts et al. 1998) and already described in Eq. (7.11); and
- **(d)** Soil water transfer dependent on a soil water potential including the effects of frozen water (e.g. Lundin 1989), represented in a proxy way by Eq. (7.66).

The latent-heat effects are described in the following. The main impact will be to delay the soil cooling in the beginning of the cold period, and to delay the soil warming in spring, although the latter effect is less important because it occurs when the solar forcing is significant. Both effects make the soil temperatures less responsive to the
atmospheric forcing and damp the amplitude of the annual soil temperature cycle. More details on the soil-freezing scheme and its impact on forecasts and the model climate are described in Viterbo et al. 1999.

The soil energy equation, Eq. (7.45), is modified in the presence of soil water phase changes as

\[
(pC)_\text{soil} \frac{dT}{dt} = \frac{\partial}{\partial z} \left[ \lambda \frac{\partial T}{\partial z} \right] + L_{\text{fus}} \rho_w \frac{\partial \theta_1}{\partial t} \tag{7.51}
\]

where \( \theta_1 \) is the volumetric ice-water content. Without loss of generality, for the grid squares characteristic of NWP models it can be assumed that

\[
\theta_1 = \theta_1(\theta, T) = f(T)\theta \tag{7.52}
\]

where \( \theta \) is the total soil-water content (liquid + ice), and

\[
f_\theta(T) = \begin{cases} 
0 & T > T_{f1} \\
0 < f_\theta(T) < 1 & T_{f1} \leq T \leq T_{f2} \\
1 & T < T_{f2}
\end{cases} \tag{7.53}
\]

where \( T_{f1} \) and \( T_{f2} \) are characteristic temperatures limiting the phase change regime. In reality, the values of \( T_{f1} \) and \( T_{f2} \) and the function \( f_\theta(T) \) have complicated dependencies on soil texture and composition (see e.g. Williams and Smith 1989), but here they are approximated in a simple way. For an idealized homogeneous, one-component soil, \( f_\theta(T) \) would be a step-function. The physical reasons for having an interval over which melting/freezing is active, rather than a threshold temperature, include (Williams and Smith 1989):

- (a) Adsorption, resulting from forces between the mineral parts of the soil and the water;
- (b) Capillarity, related to the fact that the water-free surface is not plane;
- (c) Depression of the freezing point due to the effect of dissolved salts; and
- (d) Soil heterogeneity.

To avoid an undesirable coupling between the temperature and water equations in the soil, Eq. (7.52) is simplified to

\[
\theta_1 = f_\theta(T)\theta_1 \tag{7.54}
\]

where \( \theta_1 \) is a constant, representing the amount of soil water that can be frozen (thawed). For simplicity, \( \theta_1 = (c_{\text{f1}} + c_{\text{f2}})\theta_{\text{cap}} \). The scaling with the vegetated fractions is the simplest way of distinguishing between dry (vegetation-sparse areas, e.g. deserts) and wet (vegetated) areas. Combining Eq. (7.54) with Eq. (7.51) results in

\[
\left[ (pC)_\text{soil} - L_{\text{fus}} \rho_w \frac{\partial f_\theta}{\partial T} \right] \frac{\partial T}{\partial t} = \frac{\partial}{\partial z} \left[ \lambda \frac{\partial T}{\partial z} \right] \tag{7.55}
\]

showing that the effect of freezing can be interpreted as an additional soil heat capacity, sometimes referred in the literature as the ‘heat-capacity barrier’ around freezing; not considering the process of soil water freezing/melting can lead to very large artificial temperature changes that do not occur in nature when sufficient soil water is available.

Finally, function \( f_\theta(T) \), is given by
\[ f_{11}(T) = \begin{cases} 0 & T > T_{f1} \\ 0.5 \left( 1 - \sin \left( \frac{\pi (T - 0.5 T_{f1} - 0.5 T_{f2})}{T_{f1} - T_{f2}} \right) \right) & T_{f2} \leq T \leq T_{f1} \\ 1 & T < T_{f2} \end{cases} \] (7.56)

with \( T_{f1} = T_0 + 1 \), \( T_{f2} = T_0 - 3 \).

### 7.6 Soil-Water Budget

The vertical movement of water in the unsaturated zone of the soil matrix obeys the following equation (see Richards (1931), Philip (1957), Hillel (1982), and Milly (1982) for the conditions under which Eqs. (7.57) and (7.58) are valid) for the volumetric water content \( \theta \) :

\[
\rho_w \frac{\partial \theta}{\partial t} = - \frac{\partial F_w}{\partial z} + \rho_w S_\theta
\] (7.57)

\( \rho_w \) is the water density (kg m\(^{-3}\)), \( F_w \) is the water flux in the soil (positive downwards, kg m\(^{-2}\) s\(^{-1}\)), and \( S_\theta \) is a volumetric sink term (m\(^3\) m\(^{-3}\) s\(^{-1}\)), corresponding to root extraction. Using Darcy’s law, \( F_w \) can be specified as:

\[
F_w = \rho_w \left( \lambda \frac{\partial \theta}{\partial z} - \gamma \right)
\] (7.58)

\( \lambda \) (m\(^2\) s\(^{-1}\)) and \( \gamma \) (m s\(^{-1}\)) are the hydraulic diffusivity and hydraulic conductivity, respectively.

Replacing (7.58) in (7.57), specifying \( S_\theta = S_\theta(\theta, z) \), and defining parametric relations for \( \lambda \) and \( \gamma \) as functions of soil water, a partial differential equation for \( \theta \) is obtained; it can be numerically integrated if the top boundary condition is precipitation minus evaporation minus surface runoff. The bottom boundary condition assumes free drainage. Abramopoulos et al. (1988) specified free drainage or no drainage, depending on a comparison of a specified geographical distribution of bedrock depth, with a model-derived water-table depth. For the sake of simplicity the assumption of no bedrock everywhere has been adopted.

#### 7.6.1 Interception

The interception reservoir is a thin layer on top of the soil/vegetation, collecting liquid water by the interception of rain and the collection of dew, and evaporating at the potential rate. The water in the interception reservoir, \( W_1 \), obeys

\[
\rho_w \frac{\partial W_1}{\partial t} = c_i E_i + D + I
\] (7.59)

where \( c_i E_i \) is the water evaporated by the interception reservoir (or dew collection, depending on its sign), \( D \) represents the dew deposition from other tiles, and \( I \) (kg m\(^{-2}\) s\(^{-1}\)) is the interception—the fraction of precipitation that is collected by the interception reservoir and is later available for potential evaporation. Because the interception reservoir has a very small capacity (a maximum of the order of 1 mm, see Eq. (7.2)), it can fill up or evaporate completely in one time step; special care has to be taken in order to avoid numerical problems when integrating Eq. (7.59). In addition, since \( E_i \) is defined in the vertical diffusion code, it might impose a rate of
evaporation that depletes entirely the interception layer in one time step. In order to conserve water in the atmosphere-intercepted water–soil continuum, the mismatch of evaporation of tile 3 plus dew deposition from the other tiles (which is not explicitly dealt with by the vertical diffusion) as seen by the vertical diffusion and the intercepted water has to be fed into the soil.

The equation is solved in three fractional steps: evaporation, dew deposition, and rainfall interception. The solver provides as outputs

(a) the interception layer contents at time step \( n + 1 \), \( W_i^{n+1} \);
(b) Throughfall (ie, rainfall minus intercepted water); and
(c) The evaporation effectively seen by the intercepted layer in each tile \( i \).

First, the upward evaporation \((E_i < 0)\) contribution is considered; because \( c_i E_i \) depends linearly on \( W_i \) (see Eq. (7.2)), an implicit version of the evaporating part of (7.59) is obtained by linearizing \( c_i(W_i)E_i \):

\[
\rho_w \frac{W_i^* - W_i^l}{\Delta t} = c_i(W_i)E_i + \frac{E_{im}}{W_{im}}(W_i^* - W_i^l)
\]

(7.60)

where \( W_i^* \) is the new value of interception-reservoir content after the evaporation process has been taken into account. After solving for \( W_i^* \), a non-negative value of evaporation is obtained and the evaporation seen by this fractional time step is calculated

\[
W_i^l = \max(0, W_i^*)
\]

(7.61)

\[
E_i^l = \rho_w \frac{W_i^l - W_i^l}{\Delta t}
\]

The dew deposition is dealt with explicitly for each non-snow tile in succession, for tiles 3, 4, 6, 7, 8, where tile 7 is also considered because in the exposed snow tile, the canopy is in direct evaporative contact with the atmosphere. When the evaporative flux is downwards \((E_i > 0)\)

\[
W_i^2 = W_i^l + \min\left(W_{im} - W_i^l, \frac{\Delta t}{\rho_w} c_i D_i\right)
\]

(7.62)

\[
D_i = \rho_w \frac{W_{im}^2 - W_i^l}{\Delta t}
\]

where superscript 2 denotes the final value at the end of the this fractional time step.

The interception of rainfall is considered by applying the following set of equations to large-scale and convective rainfall

\[
W_i^3 = W_i^2 + \min\left(W_{im} - W_i^2, \frac{\Delta t}{\rho_w} b_i (c_{H} + c_{L}) R_{ls}\right)
\]

\[
T_{ls} = R_{ls} - \rho_w \frac{W_i^3 - W_i^2}{\Delta t}
\]

(7.63)

\[
W_i^{*+1} = W_i^3 + \min\left(W_{im} - W_i^3, \frac{\Delta t}{\rho_w} b_i (c_{H} + c_{L}) \frac{R_{cv}}{F_{cv}}\right)
\]

\[
T_{cv} = R_{cv} - \rho_w \frac{W_i^{*+1} - W_i^3}{\Delta t}
\]
$R_{cv}/F_{cv}$ is a modified convective rainfall flux, computed by applying the heterogeneity assumption that convective rainfall only covers a fraction $F_{cv} = 0.5$ of the grid box, $b_I = 0.25$ is a coefficient of efficiency of interception of rain. The total evaporation seen by the interception reservoir is $D_I$ for tiles 4, 6, 7, and 8 and $c_I E_I + D_I$ for tile 3.

The interception reservoir model described in this section is probably the simplest water-conserving formulation based on Rutter's original proposition (Rutter et al. 1972; Rutter et al. 1975). For more complicated formulations still based on the Rutter concept see, for instance, Mahfouf and Jacquemin (1989), Dolman and Gregory (1992), and de Riddler (2001).

### 7.6.2 Soil properties

Integration of Eqs. (7.57) and (7.58) requires the specification of hydraulic conductivity and diffusivity as a function of soil-water content. Mahrt and Pan (1984) have compared several formulations for different soil types. The widely used parametric relations of Clapp and Hornberger (1978) (see also Cosby et al. 1984) are adopted:

$$\gamma = \gamma_{sat}\left(\frac{\theta}{\theta_{sat}}\right)^{2b + 3}$$

$$\lambda = \frac{b \gamma_{sat}(\psi_{sat})}{\theta_{sat}} \left(\frac{\theta}{\theta_{sat}}\right)^{b + 2}$$

(7.64)

$b$ is a non-dimensional exponent, $\gamma_{sat}$ and $\psi_{sat}$ are the values of the hydraulic conductivity and matric potential at saturation, respectively. A minimum value is assumed for $\lambda$ and $\gamma$ corresponding to permanent wilting-point water content.

Cosby et al. (1984) tabulate best estimates of $b$, $\gamma_{sat}$, $\psi_{sat}$ and $\theta_{sat}$, for the 11 soil classes of the US Department of Agriculture (USDA) soil classification, based on measurements over large samples. Since the model described here specifies only one soil type everywhere, and because the determination of the above constants is not independent of the values of $\theta_{cap}$ and $\theta_{pw}$, the following procedure is adopted.

A comprehensive review of measurements of $\theta_{cap}$ and $\theta_{pw}$ may be found in Patterson (1990). Starting from Patterson's estimates of $\theta_{cap}$ and $\theta_{pw}$ for the 11 USDA classes, a mean of the numbers corresponding to the medium-texture soils (classes 4, 5, 7, and 8, corresponding to silt loam, loam, silty clay loam and clay loam, respectively) is taken. The resulting numbers are $\theta_{cap} = 0.323$ $m^3$ $m^{-3}$ and $\theta_{pw} = 0.171$ $m^3$ $m^{-3}$. Averaging the values of Cosby et al. (1984) for soil moisture and soil-water conductivity at saturation for the same classes gives the numerical values $\gamma_{sat} = 5.57 \times 10^{-6}$ $m$ $s^{-1}$ and $\theta_{sat} = 0.472$ $m^3$ $m^{-3}$. The Clapp and Hornberger expression for the matric potential

$$\psi = \psi_{sat}\left(\frac{\theta}{\theta_{sat}}\right)^{b}$$

(7.65)

is used with $\psi(\theta_{pw}) = -153$ $m$ ($-15$ bar) and $\psi(\theta_{cap}) = -3.37$ $m$ ($-0.33$ bar) (see Hillel 1982; Jacquemin and Noilhan 1990) to find the remaining constants $b$ and $\psi_{sat}$. The results are $b = 6.04$ and $\psi_{sat} = -0.338$ $m$.

The above process ensures a soil that has an availability corresponding to the average value of medium-texture soils, and yields a quantitative definite hydraulic meaning to $\theta_{cap}$ and $\theta_{pw}$ compatible with the Clapp and Hornberger relations (see Table 7.2 for a summary of the soil constants).

Finally, the water transport in frozen soil is limited in the case of a partially frozen soil, by considering the effective hydraulic conductivity and diffusivity to be a weighted average of the values for total soil water and a very small
value (for convenience, taken as the value of Eq. (7.64) at the permanent wilting point) for frozen water. The soil properties, as defined above, also imply a maximum infiltration rate at the surface defined by the maximum downward diffusion from a saturated surface. If the throughfall exceeds the maximum infiltration rate, the excess precipitation is put into runoff. However, in practice the maximum infiltration rate is so large that this condition is never reached. Surface runoff is therefore only produced if the soil becomes saturated.

7.6.3 Discretization and the root profile

A common soil discretization is chosen for the thermal and water soil balance for ease of interpretation of the results, proper accounting of the energy involved in freezing/melting soil water, and simplicity of the code. Equations Eqs. (7.57) and (7.58) are discretized in space in a similar way to the temperature equations, i.e., soil water and root extraction defined at full layers, $\theta_k$ and $\rho_{\text{water}} S_{0,k}$, and $F_{k+1/2}$ the flux of water at the interface between layer $k$ and $k + 1$. The resulting system of equations represents an implicit, water-conserving method.

For improved accuracy, the hydraulic diffusivity and conductivity are taken as (see Mahrt and Pan 1984)

$$\lambda_{k+1/2} = (1 - f_{tr}^*)\lambda [\max(\theta_k^n, \theta_{k+1}^n)] + f_{tr}^* \lambda(\theta_{\text{pwp}})$$  \hspace{1cm} (7.66)

$$\gamma_{k+1/2} = (1 - f_{tr}^*)\gamma [\max(\theta_k^n, \theta_{k+1}^n)] + f_{tr}^* \gamma(\theta_{\text{pwp}})$$

where $f_{tr}^* = \min[f_{tr}(\theta_k), f_{tr}(\theta_{k+1})]$. The boundary conditions are given by

$$F_{1/2} = \rho_w \gamma_{1/2}$$ \hspace{1cm} (7.67)

The difference between throughfall $T$ and surface runoff $Y_{\text{sf}}$ (kg m$^{-2}$ s$^{-1}$) is the soil infiltration at the surface:

$$T = T_{is} + T_{cv}$$

$$Y_{\text{sf}} = \max(0, T_{is} + M_{sn} - I_{f, mx}) + \frac{\max(0, F_{cv} T_{cv} - I_{f, mx})}{F_{cv}}$$ \hspace{1cm} (7.68)

$$I_{f, mx} = \rho_w \left[ \frac{\lambda_{1/2} (\theta_{\text{sat}} - \theta_i)}{0.5D_{3j}} + \gamma_{1/2} \right]$$

and $\lambda_{1/2} = f_{tr}^* \lambda(\theta_{\text{pwp}}) + (1 - f_{tr}^*) \lambda(\theta_{\text{sat}})$, with a similar equation for $\gamma_{1/2}$. The evaporation at the top of the soil layer, $E_{1/2}$, is computed as the sum of the evaporations of tile 8 plus the contributions necessary to conserve water with the solver of the interception layer:

(a) tile 3 mismatch(after the evaporated water used by the interception reservoir for the given tile is subtracted); and

(b) when the evaporative fluxes are downward (i.e., dew deposition), the evaporation for tiles 4, 6 and the canopy evaporation of tile 7.

Root extraction is computed as

$$\rho_w S_{0,k} = \sum_i C_i R_i \theta_k / \sum_j R_j \theta_j$$ \hspace{1cm} (7.69)
where the sum over tiles \( i \) is done for tiles 4, 6, and 7 (for which only the transpiration is used) and the sum over \( j \) is done over all soil levels. In case of dew deposition (i.e., tile downward evaporative flux), \( S_{0,h} = 0 \).

### 7.7 Sea/Lake Ice

Any non-land point (i.e., a grid point with land cover less or equal 0.5) can have two fractions, open water and ice. A surface analysis defines the ice fraction, \( c_i \), and the temperature of the open water fraction; both quantities are kept constant during the forecast. No distinction is made between surface and skin temperature for the open water fraction (see Table 7.2).

The ice fraction is modelled as an ice slab, with open water underneath and a skin temperature for the thermal contact with the atmosphere. The main caveats in the sea ice parameterization are:

(a) Fixed depth of the slab (which can be relaxed once there is a reliable data set to specify its geographic distribution;

(b) Fixed fraction, which is a reasonable assumption for a 10-day forecast period, and avoids the need for the momentum balance of the ice and its complex rheology (see, e.g., Flato and Hibler 1992) and the definition of the ocean currents;

(c) No snow accumulation on top of the ice (although one of the main effects of snow, i.e., a markedly different surface albedo, is partially emulated by the prescribed seasonal albedo in Table 2.2).

The ice heat transfer is assumed to obey the following Fourier law of diffusion

\[
(\rho C)_I \frac{\partial T_I}{\partial t} = \frac{\partial}{\partial z} \left[ \lambda_I \frac{\partial T_I}{\partial z} \right]
\]  
(7.70)

where \((\rho C)_I = 1.88 \times 10^6 \text{J m}^{-3} \text{K}^{-1}\) is the volumetric ice heat capacity, \(T_I\) is the ice temperature, and \(\lambda_I = 2.03 \text{ W m}^{-1} \text{K}^{-1}\) is the ice thermal conductivity. The boundary condition at the bottom is the temperature of the frozen water, \(T_n = T_0 - 1.7\) and the top boundary condition is the net heat flux at the surface, obtained from the solution of the ice skin thermal budget.

Eq. (7.70) is solved with the ice discretized in four layers, with the depth of the top three layers as in the soil model and the depth of the bottom layer defined as

\[
D_{I,4} = D_I - \sum_{j=1}^{3} D_{I,j}
\]  
(7.71)

and the total depth of the ice slab, \(D_I\), is prescribed as 1.5 m. In order to ensure a constant ice fraction, the solution of the ice thermal budget is capped to the ice melting temperature, \(T_m = T_0\) at all levels. The details of the numerical discretization can be found in Section 7.8.

### 7.8 Numerical Solution of the Surface Equations

#### 7.8.1 Recap of the analytical equations

The water budget (Eqs. (7.57)–(7.58), with boundary conditions given by Eq. (7.67)), the soil energy budget (Eq. (7.45), with boundary conditions given by Eqs. (7.46)–(7.47)) and the ice energy budget (Eq. (7.70)) can be rewritten in a generalised form as:
\[
\frac{\partial \Psi}{\partial t} = \frac{1}{C} \frac{\partial}{\partial z} \left( \lambda \frac{\partial \Psi}{\partial z} - \gamma \right) + S_\Psi
\]

(7.72)

The meaning of the different variables in each individual equations is summarized Table 7.6, together with the respective upper and lower boundary conditions, \( F_\Psi \).

<table>
<thead>
<tr>
<th>Equation</th>
<th>( \Psi )</th>
<th>( C )</th>
<th>( \lambda )</th>
<th>( \gamma )</th>
<th>( S_\Psi )</th>
<th>UBC</th>
<th>LBC</th>
</tr>
</thead>
<tbody>
<tr>
<td>Soil moisture</td>
<td>0</td>
<td>1</td>
<td>( \lambda_0 )</td>
<td>( \gamma_0 )</td>
<td>( S_0 )</td>
<td>( F_0 = 1 f - c_k E_N )</td>
<td>( F_0 = \gamma_0 )</td>
</tr>
<tr>
<td>Soil temperature</td>
<td>( T )</td>
<td>( (\rho C)_e )</td>
<td>( \lambda_T )</td>
<td>0</td>
<td>0</td>
<td>( F_T = H^N )</td>
<td>( F_T = 0 )</td>
</tr>
<tr>
<td>Ice temperature</td>
<td>( T_I )</td>
<td>( (\rho C)_I )</td>
<td>( \lambda_I )</td>
<td>0</td>
<td>0</td>
<td>( F_I = H^N )</td>
<td>( T_{N+1} = T_{0,I} )</td>
</tr>
</tbody>
</table>

UBC and LBC stand for upper and lower boundary condition, respectively.

7.8.2 Implicit numerical solution

Eq. (7.72) is time discretized in the following way:

\[
\frac{\Psi^{t+1} - \Psi^t}{\Delta t} = \frac{1}{C} \frac{\partial}{\partial z} \left( \lambda \frac{\partial \Psi}{\partial z} - \gamma \right) + S_\Psi
\]

(7.73)

where

\[
\Psi^{t+1} = \alpha_{\text{impl}} \Psi^{t+1} + (1 - \alpha_{\text{impl}}) \Psi^t
\]

(7.74)

and the semi-implicit coefficient, \( \alpha_{\text{impl}} = 1 \). If the prognostic variable \( \Psi \) is defined at full levels and the fluxes \( F_\Psi \) are defined at half-levels (the interface between layers), Eq. (7.73) can be discretized in space to give:

\[
\frac{\Psi - \Psi^t}{\alpha_{\text{impl}}} = \frac{\Delta t}{C_k} \left( \frac{\lambda_{k-1/2} (\Psi_{k-1/2} - \Psi_{k}) - \lambda_{k-1/2} (\Psi_{k} - \Psi_{k+1})}{\Delta z_k \Delta z_{k-1/2}} + \frac{\gamma_{k-1/2} - \gamma_{k+1/2}}{\Delta z_k} \right) + \Delta t S_{\Psi,k} \quad k = 2, \ldots, N\delta - 1
\]

(7.75)

\[
\frac{\Psi - \Psi^t}{\alpha_{\text{impl}}} = \frac{\Delta t}{C_k} \left( \frac{F_{\Psi}^T \left( \Psi_{k} - \Psi_{k+1} \right) - \lambda_{k-1/2} (\Psi_{k} - \Psi_{k+1})}{\Delta z_k \Delta z_{k-1/2}} + \frac{\gamma_{k-1/2} - \gamma_{k+1/2}}{\Delta z_k} \right) + \Delta t S_{\Psi,k} \quad k = 1
\]

(7.75)

\[
\frac{\Psi - \Psi^t}{\alpha_{\text{impl}}} = \frac{\Delta t}{C_k} \left( \frac{\lambda_{k-1/2} (\Psi_{k-1/2} - \Psi_{k}) - \lambda_{k-1/2} (\Psi_{k} - \Psi_{k+1})}{\Delta z_k \Delta z_{k-1/2}} + \frac{\gamma_{k-1/2} - \gamma_{k+1/2}}{\Delta z_k} \right) + \Delta t S_{\Psi,k} \quad k = N\delta
\]

(7.75)

where the horizontal brace means that the term exists only for the ice temperature equation (because of the bottom temperature boundary condition for ice) and \( \Delta z_k, \Delta z_{k-1/2} \) and \( \Delta z_{k+1/2} \) represent the thickness of layer \( k \), and the depths of its centre, the top and the bottom interface, respectively:

\[
\Delta z_k = D_k = z_{k+1/2} - z_{k-1/2}
\]

\[
\Delta z_{k+1/2} = z_{k+1} - z_k
\]

(7.76)
Eq. (7.75) leads to a tri-diagonal system of equations:

\[
\frac{\hat{\Psi}_{k-1}}{\alpha_{impl}} \left( \hat{\lambda}_{k-1/2} \right) + \frac{\hat{\Psi}_{k}}{\alpha_{impl}} \left( 1 + \hat{\lambda}_{k+1/2} \right) - \frac{\hat{\Psi}_{k+1}}{\alpha_{impl}} \left( \hat{\lambda}_{k+1/2} \right) = \frac{\Psi_{impl}'}{\Delta t} \left( \gamma_{k-1/2} - \gamma_{k+1/2} \right) + \Delta t S_{\Psi,k} \quad k = 2, \ldots, Ns - 1
\]

\[
\frac{\hat{\Psi}_{k}}{\alpha_{impl}} \left( 1 + \hat{\lambda}_{k+1/2} \right) - \frac{\hat{\Psi}_{k+1}}{\alpha_{impl}} \left( \hat{\lambda}_{k+1/2} \right) = \frac{\Psi_{impl}'}{\Delta t} \left( \gamma_{k-1/2} - \gamma_{k+1/2} \right) + \Delta t S_{\Psi,k} \quad k = 1
\]

\[
\frac{\hat{\Psi}_{k-1}}{\alpha_{impl}} \left( \hat{\lambda}_{k-1/2} \right) + \frac{\hat{\Psi}_{k}}{\alpha_{impl}} \left( 1 + \hat{\lambda}_{k+1/2} \right) - \frac{\hat{\Psi}_{k+1}}{\alpha_{impl}} \left( \hat{\lambda}_{k+1/2} \right) = \frac{\Psi_{impl}'}{\Delta t} \left( \gamma_{k-1/2} - \gamma_{k+1/2} \right) + \Delta t S_{\Psi,k} \quad k = Ns
\]

with the generalized modified diffusivities, \( \hat{\lambda}_{k-1/2} \), defined as:

\[
\hat{\lambda}_{k-1/2} = \frac{\Delta t \alpha_{impl} \lambda_{k-1/2}}{\Delta z_{k-1/2}}
\]

\[
I \begin{bmatrix} \Delta z_{Ns} \gamma_{k-1/2} \\ \Psi_{Ns+1} \end{bmatrix} = D_{Ns/2}
\]

\[
\Psi_{Ns+1} = T_{0,I}
\]

where \( D_{Ns} \) is the depth of the deepest soil layer. The discretization above conserves water (energy) and is linearly stable. The coefficients \( \lambda \) and \( \gamma \) are a function of variable at the current time step, \( \Psi^\alpha \).

### 7.9 Code

The surface parameterization computations are shared between the vertical diffusion routine (VDFMAIN, see Chapter 3) and the main surface routine, SRFMAIN. In VDFMAIN, the tile fluxes and skin temperatures are computed: After the elimination part of the tri-diagonal system of equations is computed, the energy budget for each tile is computed before back-substitution.

At the start of the model integration, the following setup routine is called to initialize a module specific to the soil-code:

- **SUSOIL.** Setup routine for soil/snow/ice constants.

The main subroutine of the surface code (SRFMAIN) is called from CALLPAR, with: (a) values of the surface prognostic equations at time step n, convective and large scale rainfall and snowfall, tile evaporation, sensible and latent heat fluxes, and temperatures, net surface longwave flux, tile net shortwave flux as inputs; and (b) tendencies for the surface prognostic variables, plus a comprehensive set of diagnostic arrays as outputs. SRFMAIN does a sequence of computations and subroutine calls:
Part IV: ‘Physical processes’

- **SRFSN.** Solution of the snow energy and water budget and computation of the next time step density and albedo fields. Inputs: snow depth, temperature, density and albedo at the current time step, soil temperature, shortwave and longwave radiation fluxes, snowfall, and tile fluxes. Outputs: snow depth, temperature, density and albedo at the next time step, meltwater flux, and basal heat flux.

- **SRFRCG.** Computes apparent soil heat capacity, i.e., including effects of soil freezing. Inputs: soil temperature and vegetation covers. Output is volumetric heat capacity.

- **SRFT.** Solution of the soil heat budget. Inputs: Soil temperature, soil moisture, longwave radiative flux, snow basal heat flux, volumetric heat capacity, tile evaporation, sensible heat flux and shortwave radiative flux. Output: Soil temperature at the next time step. First the modified heat diffusivity, the soil energy per unit area and the right-hand side of the system of equations are computed. The generalized surface tridiagonal solver, **SRFWDIF,** is called to solve for the semi-implicit variable, \( \hat{T}/\alpha \). The soil temperatures for the next time step are computed at the end.

- **SRFI.** Solution of the ice heat budget. Inputs: Ice temperature, longwave radiative flux, tile evaporation, sensible heat flux and shortwave radiative flux. Output: Ice temperature at the next time step. First the modified heat diffusivity, the ice energy per unit area and the right-hand side of the system of equations are computed. The generalized surface tridiagonal solver, **SRFWDIF,** is called to solve for the semi-implicit variable, \( \hat{T}/\alpha \). The ice temperatures for the next time step are computed at the end.

- **SRFWL.** Solution of the interception layer water budget. Inputs: Interception layer contents, low and high vegetation water cover, maximum capacity of the interception layer, convective and large scale rainfall, snow evaporation of shaded snow tile, and tile evaporation. Outputs: Interception layer at next time step, convective and large scale throughfall and tile evaporation collected (or depleting) the interception layer.

- **SRFWEXC.** First part of the computation of the soil water budget, i.e., computation of the coefficients of the tridiagonal system of equations for \( \hat{\Theta} \). This includes the partitioning of transpiration into root extraction at the different layers and soil hydraulic coefficients including the effect of frozen water. Inputs: Soil moisture and temperature, convective and large-scale throughfall, snowmelt, tile evaporation, tile evaporation collected (or depleting) the interception layer, and snow evaporation of the shaded snow tile. Outputs: Modified diffusivity for water, right-hand side of the tridiagonal system, and layer depths.

- **SRFWDIF.** Generalized surface tridiagonal solver. Inputs: Values of \( \psi \) at the current time step, generalized modified diffusivities, soil energy (or water) per unit area, and right-hand side of equations. Output: \( \psi/\alpha \). The routine computes the coefficients on the left-hand side of the equations and solves the equations using and LU-decomposition and back substitution in one downward scan and one upward scan.

- **SRFWINC.** Computation of next time step soil water. Inputs: \( \hat{\Theta}/\alpha \) and current time step soil water. Output: next time step soil water.

- **SRFWNG.** Bounded-value operator for intercepted water (limited to non-negative values and values below or equal the maximum contents of the interception layer) and soil water (limited to non-negative values and values below or equal saturation). The “soil column” is scanned from top to bottom and the amount of water needed to satisfy physical limits in each layer are borrowed from the layer below. The water exchanged in this way is accounted for as runoff. Inputs: next time step intercepted water and soil water. Output: Bounded values of the same quantities.

Relevant routines from the vertical diffusion code, discussed in full detail in Chapter 3, include:

- **SUVEG.** Assignment of vegetation related constants.
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- **VDFBC.** Definition of tile fractions and related characteristics.
- **VDFSURF.** Definition of bare soil resistance, low and high canopy resistances.
- **VDFEXCS.** Computation of aerodynamical part of exchange coefficients for heat and moisture, including stability computations.
- **VDFEVAP.** Computation of evapotranspiration for each tile.
- **VDFSFLX.** Surface fluxes for each tile, defined at time $t$.
- **VDFTSK.** Computation of the tile skin temperatures, as a the solution of the tile energy balance.
- **VDFTFLX.** Computation of the tile fluxes at time $t + 1$. 
Part IV: PHYSICAL PROCESSES

CHAPTER 8 Methane oxidation

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8.2 Methane oxidation
8.3 The parametrization
  8.3.1 Methane oxidation
  8.3.2 Photolysis in the mesosphere
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8.1 INTRODUCTION

A study of stratospheric humidity in analyses and multi-year simulations has shown that the ECMWF system prior to 1999 was capable of producing a broadly realistic distribution of water vapour at, and immediately above, the tropopause, and that the slow upward transfer of water vapour in the tropical stratosphere could be captured quite reasonably given sufficiently fine vertical resolution in the model (Simmons et al. 1999). However, values of water vapour in the tropical upper stratosphere, and throughout much of the extratropical stratosphere, were too low. This deficiency has now been remedied by the introduction of a simple parametrization of the upper-stratospheric moisture source due to methane oxidation. A sink representing photolysis in the mesosphere is also included. The scheme was derived as a simplification of an approach adopted by Peter Stott and Anne Pardaens at the Department of Meteorology, University of Edinburgh, notes on which and helpful references were supplied by Bob Harwood.

8.2 METHANE OXIDATION

Methane is produced by natural and anthropogenic sources at the earth’s surface, and is well-mixed in the troposphere. Its volume mixing ratio is currently around 1.7 ppmv. It is carried upwards in the tropical stratosphere and decreases in relative density (due to oxidation) to values of around 0.2–0.4 ppmv around the stratopause. Mean stratospheric descent at higher latitudes results in relatively low values of methane at these latitudes in the middle and lower stratosphere.

Brasseur and Solomon (1984) provide an account of the chemistry of carbon compounds in the stratosphere and mesosphere. The long chain of reactions starting from methane (CH₄) ends with the production of water vapour (H₂O) and molecular hydrogen (H₂) in the stratosphere and mesosphere. This occurs such that the sum

\[ 2[CH₄] + [H₂O] + [H₂] \]

is approximately uniformly distributed in the absence of precipitation, where \([\_\_]\) denotes a volume mixing ratio. Le Texier et al. (1988) provide calculations of the relative amounts of H₂O and H₂, showing that the predominant production is that of water vapour in the vicinity of the stratopause. They indicate, however, that H₂ production in the mesosphere, and relatively strong descent in winter and early spring at high latitudes, may result in the upper
stratosphere being relatively dry in these seasons and latitudes.

There is, nevertheless, good observational evidence that over much of the stratosphere the quantity

$$2[CH_4] + [H_2O]$$

is quite uniformly distributed with a value somewhat over 6 ppmv. Jones et al. (1986) provide evidence for this from the LIMS and SAMS instruments on the Nimbus 7 satellite launched in 1978, and a particularly clear demonstration is given by Bithell et al. (1994) based on HALOE data from the UARS satellite. In a pressure–latitude section at about the austral spring equinox, Bithell et al. show the result to fail significantly only below 10 hPa in the high-latitude southern hemisphere due, presumably, to condensation at the very cold temperatures in the Antarctic polar vortex.

Prior to cycle 25r1 of the IFS, the parametrization used the value 6 ppmv for the sum $2[CH_4] + [H_2O]$. This version was used in production of the ERA-40 reanalyses, which have been found to be generally drier in the stratosphere than the climatology derived by Randel et al. (1998) from UARS measurements. From cycle 25r1 onwards, the parametrization uses the value 6.8 ppmv, based on Randel et al.’s data as presented in Fig. 8.1.

Figure 8.1 Annual-mean distribution of the sum of twice the volume mixing ratio of methane and of the mixing ratio of water vapour (ppmv) as a function of pressure and potential vorticity (expressed as equivalent latitude), derived from UARS (HALOE, supplemented by CLAES and MLS) data analysed by Randel et al. (1998). The contour interval is 0.1 ppmv, and shading denotes the range 6.6-6.9 ppmv.
8.3 THE PARAMETRIZATION

8.3.1 Methane oxidation
We assume that the volume mixing ratio of water vapour $[H_2O]$ increases at a rate

$$2k_1[CH_4]$$  \hspace{1cm} (8.1)

We further assume that

$$2[CH_4] = 6.8 \text{ ppmv} - [H_2O]$$  \hspace{1cm} (8.2)

The rate of increase of volume mixing ratio of water vapour (in ppmv) is thus

$$k_1(6.8 - [H_2O])$$  \hspace{1cm} (8.3)

In terms of specific humidity, $q$, the source is

$$k_1(Q - q)$$  \hspace{1cm} (8.4)

where (having divided by $1.6 \times 10^6$ to convert from volume mixing ratio in ppmv to specific humidity) the parameter $Q$ has the value $4.25 \times 10^{-6}$, or 4.25 mg/kg.

The rate $k_1$ could be determined, for example, from a 2-D model with comprehensive chemistry, as in the scheme developed at Edinburgh University. However, in this first scheme for use at ECMWF we prescribe a simple analytical form for $k_1$ which varies only with pressure.

The photochemical life time of water vapour is of the order of 100 days near the stratopause, 2000 days at 10 hPa, and effectively infinite at the tropopause (Brasseur and Solomon 1984). A prescription of $k_1$ that gives a reasonable profile up to the stratopause is provided by

$$k_1 = \frac{1}{86400 \tau_1}$$  \hspace{1cm} (8.5)

where $k_1$ is given in s$^{-1}$ and the timescale, $\tau_1$, in days, is given in terms of pressure, $p$, in Pa, by:

$$\tau_1 = \begin{cases} 100 & p \leq 50 \\ 100 \left[1 + \alpha_1 \left(\frac{\ln(p/50)}{\ln(10000/p)}\right)^4\right] & 50 < p < 10000 \\ \infty & p \geq 10000 \end{cases}$$  \hspace{1cm} (8.6)

where we define

$$\alpha_1 = \frac{19\ln 10}{(\ln 20)^3}$$  \hspace{1cm} (8.7)

to give a time-scale of 2000 days at the 10 hPa level.
This parametrization moistens rising air in the tropical stratosphere. This air will earlier have been freeze-dried near the tropopause, where specific humidities can locally fall well below 1 mg/kg. Specific humidities approaching the value $Q$ will be reached near the stratopause. Descent near the poles will bring down air with specific humidity close to $Q$. Expression (8.4) will then yield a source term that is weaker in polar than in tropical latitudes, so reasonable results may be obtained without imposing a latitudinal variation of $k_1$. (Strictly, $k_1$ should vanish in the polar night, where photodissociation does not produce the excited oxygen $O(^3D)$, which in turn produces the OH radical, these two species being intimately involved in the production of water vapour from methane).

### 8.3.2 Photolysis in the mesosphere

For model versions with an uppermost level at 0.1 hPa, or lower, there is no strong need to include the sink of water vapour that occurs in the mesosphere and above due to photolysis. However, for completeness we include a simple representation of this effect, modifying the source term (8.4) by adding a decay term $-k_2 q$ above a height of about 60 km. The full source/sink term becomes

$$k_1 (Q - q) - k_2 q$$

(8.8)

As for $k_1$ we take $k_2$ independent of latitude with parameters chosen to match the vertical profile of photochemical lifetime presented by Brasseur and Solomon (1984). Specifically,

$$k_2 = \frac{1}{86400 \tau_2}$$

(8.9)

with

$$\tau_2 = \begin{cases} 3 & p \leq 0.1 \\ \left\{ \exp\left(\alpha_2 - 0.5(\ln 100 + \alpha_2) \left(1 + \cos \frac{\pi \ln(p/20)}{\ln 0.005}\right) \right) - 0.01 \right\}^{-1} & 0.1 < p < 20 \\ \infty & p \geq 20 \end{cases}$$

(8.10)

and

$$\alpha_2 = \ln\left(\frac{1}{3} + 0.01\right)$$

(8.11)

The vertical profile of the photochemical lifetime of the combined scheme, $(k_1 + k_2)^{-1}$, is shown below in Fig. 8.2, in which we have converted to height as a vertical coordinate assuming an isothermal atmosphere with a temperature of 240 K. Comparison of this profile with that for $H_2O$ shown in Fig. 5.21 of Brasseur and Solomon (1984) indicates reasonable agreement.
8.4 CODE

The calculations for methane oxidation and photolysis of water vapour are performed in subroutine METHOX.

This routine calculates the tendency of water vapour due to methane oxidation and due to photolysis following (8.8). The order of the calculations is as follows:

- find time-scale for methane oxidation following (8.6)
- solve first part of (8.8)
- find time-scale for water vapour photolysis following (8.10)
- solve second part of (8.8)

The setup of the constants used in METHOX is performed in SUMETHOX which is called from SUPHEC. The constants are kept in module YOEMETH. The controlling switch for the methane oxidation is LEMETHOX which is part of namelist NAEPHY.
Part IV: PHYSICAL PROCESSES

CHAPTER 9  Ozone chemistry parametrization

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9.1 Introduction
9.2 The ECMWF ozone parameterization

9.1 INTRODUCTION
Ozone is fully integrated into the ECMWF forecast model and analysis system as an additional three-dimensional model and analysis variable similar to humidity. The forecast model includes a prognostic equation for the ozone mass mixing ratio (kg/kg)

\[
\frac{dO_3}{dt} = R_{O_3}
\]

(9.1)

where \(R_{O_3}\) is a parameterization of sources and sinks of ozone. Without such a source/sink parameterization the ozone distribution would drift to unrealistic values in integrations longer than a few weeks. The source/sink parameterization must maintain a realistic ozone distribution over several years of integration, without reducing the dynamic variability of ozone. In addition, we would like the parameterization to be able to create an Antarctic ozone hole when the conditions are right.

9.2 THE ECMWF OZONE PARAMETERIZATION
The parameterization used in the ECMWF model is an updated version of Cariolle and Déqué (1986), which has been used in the ARPEGE climate model at Météo-France. This parameterization assumes that chemical changes in ozone can be described by a linear relaxation towards a photochemical equilibrium. It is mainly a stratospheric parameterization. The relaxation rates and the equilibrium values have been determined from a photochemical model, including a representation of the heterogeneous ozone hole chemistry. The updated version of the parameterization (with coefficients provided by Pascal Simon, Météo-France) is

\[
R_{O_3} = c_0 + c_1(O_3 - \overline{O_3}) + c_2(T - \overline{T}) + c_3(O_3^{\text{uparrow}} - \overline{O_3^{\text{uparrow}}}) + c_4(\text{Cl}_{\text{EQ}})^2O_3
\]

(9.2)

where

\[
O_3^{\text{uparrow}}(p) = -\int_{p_0}^{p} \frac{O_3(p')}{\overline{g}} dp'
\]

(9.3)

Here \(c_i\) are the relaxation rates and \(\overline{T}, \overline{O_3}\), and \(O_3^{\text{uparrow}}\) are photochemical equilibrium values, all functions of latitude, pressure, and month. \(\text{Cl}_{\text{EQ}}\) is the equivalent chlorine content of the stratosphere for the actual year, and is
the only parameter that varies from year to year (see Fig. 9.1). For the ECMWF model it was necessary to replace the photochemical equilibrium values for ozone with an ozone climatology (Fortuin and Langematz, 1995) derived from observations. The heterogeneous part is only turned on below a threshold temperature of 195 K.

Figure 9.1 Equivalent chlorine content of the stratosphere in ppt for the heterogeneous chemistry part of the ozone source/sink parameterization (provided by Pascal Simon, Météo-France).
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CHAPTER 10  Climatological data

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10.3 Mean orography
10.4 Land sea mask
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10.6 Parameters for gravity-wave and orographic drag schemes
10.7 Vegetation parameters
10.8 Albedo
10.9 Aerosols
10.10 Ozone

10.1 INTRODUCTION
The ECMWF model uses a series of climate fields of different origin which have different resolution and different projections. Brankovic and Van Maanen (1985) describe a set of programs (known as PREPCLIM software) to interpolate the different fields to the requested target resolution. The software handles all the target resolutions that are in use at ECMWF and either full or reduced Gaussian grids. Grid areas at the model resolution are referred to as ECMWF or model grid squares. This appendix describes the different climate fields and the procedures to derive the fields that are needed by the model.

10.2 TOPOGRAPHIC DATA
The model orography and land use fields are based on the terrain elevation data set GTOPO30 at 30” resolution (Gesch and Larson, 1998), the terrain elevation data for Greenland KMS DEM also at 30” resolution (Ekholm, 1996) and the Global Land Cover Characteristics (GLCC) data set at 1 km resolution.

The GTOPO30 data set, as used in the IFS, was completed in 1996 through a collaborative effort led by the US Geological Survey’s Data Centre (EDC, see http://edcwww.cr.usgs.gov/landdaac/gtopo30/gtopo30.html) and was derived from a variety of information sources. It contains terrain elevation above mean sea level at a resolution of 30 arc seconds with -9999 code for sea points. A lake mask is not included.

Greenland KMS DEM replaces GTOPO30 for the Greenland area, because of the better accuracy of the Greenland data.

The Global Land Cover Characteristics (GLCC) data set has been derived from 1 year of Advanced Very High Resolution Radiometer (AVHRR) data, digital elevation models, ecoregions and map data. The nominal resolution is 1 km, and the data comes on a Goode Homolosine global projection. The data base provides for each pixel a biome classification based on several of the popular classifications, including BATS, SiB and SiB2. The BATS classifica-
tion has been adopted for the IFS because it contains inland water as one of its classes.

Due to their high resolution and global coverage, these data sets are rather big and therefore difficult to handle by the standard PREPCLIM software. Therefore the original data has been converted to an intermediate resolution of 2’30” which is much easier to handle by the standard PREPCLIM software. The derived 2’30” data set contains the following fields:

- Mean elevation above mean sea level
- Land fraction
- Lake fraction
- Fractional cover for all 20 BATS biome classes (see Table 10.1)

Finally, also the original US-Navy 10’ data is still used for the subgrid orography contribution to the roughness length. It contains the average terrain height of each grid element, as well as maximum and minimum height, number and orientation of significant ridges, and percentages of water and urban areas. In future the roughness length computation will be upgraded to make optimal use of the high resolution GTOPO30 data.

<table>
<thead>
<tr>
<th>Index</th>
<th>Vegetation type</th>
<th>H/L veg</th>
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<tr>
<td>1</td>
<td>Crops, Mixed Farming</td>
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<tr>
<td>2</td>
<td>Short Grass</td>
<td>L</td>
</tr>
<tr>
<td>3</td>
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<td>4</td>
<td>Deciduous Needleleaf Trees</td>
<td>H</td>
</tr>
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<td>5</td>
<td>Deciduous Broadleaf Trees</td>
<td>H</td>
</tr>
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<td>6</td>
<td>Evergreen Broadleaf Trees</td>
<td>H</td>
</tr>
<tr>
<td>7</td>
<td>Tall Grass</td>
<td>L</td>
</tr>
<tr>
<td>8</td>
<td>Desert</td>
<td>-</td>
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<td>9</td>
<td>Tundra</td>
<td>L</td>
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<tr>
<td>10</td>
<td>Irrigated Crops</td>
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</tr>
<tr>
<td>11</td>
<td>Semidesert</td>
<td>L</td>
</tr>
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<td>Ice Caps and Glaciers</td>
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<td>17</td>
<td>Deciduous Shrubs</td>
<td>L</td>
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<tr>
<td>18</td>
<td>Mixed Forest/woodland</td>
<td>H</td>
</tr>
<tr>
<td>19</td>
<td>Interrupted Forest</td>
<td>H</td>
</tr>
<tr>
<td>20</td>
<td>Water and Land Mixtures</td>
<td>L</td>
</tr>
</tbody>
</table>
Chapter 10 ‘Climatological data’

10.3 MEAN OROGRAPHY

Orography, or geopotential height, is derived from the 2’30” data by averaging. Source and target grid are overlaid, and weighted averages are computed by considering the fractions of source grid areas that cover the target grid square.

![Orography at T511 resolution.](image)

Figure 10.1 Orography at T511 resolution.

The orography is spectrally fitted to ensure consistency in spectral space between the orography and the model resolution. Smoothing is applied in spectral space with a $\nabla^4$ operator, where damping by a factor 5 is applied to the smallest scales. This applies to all operational resolutions (T195, T319, T455, T511). Orographic ripples appear as a consequence of the spectral fitting. Fig. 10.1 shows the orography at T511 resolution.

10.4 LAND SEA MASK

Each grid point of the model is provided with a land fraction parameter, derived from the 2’30” data. The model converts this parameter into a mask where grid points that have more than 50% land are considered as land points.

10.5 ROUGHNESS LENGTHS

The model uses ‘effective’ roughness lengths for momentum ($z_{0m}$) and heat/moisture ($z_{0h}$) in the surface boundary conditions for wind and temperature/moisture respectively. These parameters represent the effect on turbulent transport of small-scale surface elements ranging from vegetation and small-scale obstacles to subgrid orography. The roughness lengths area blend of three contributions: vegetation, urbanization and subgrid orography (see Figs. 10.2 and 10.3). The following procedure is applied:

(a) The vegetation roughness length is taken from a $5^\circ \times 5^\circ$ grid from Munich University (Baumgartner et al., 1977) and interpolated to the requested model grid. Symbols $z_{0mVeg}$ and $z_{0hVeg}$ will be used for the pure vegetation roughness lengths.
(b) The fraction of urbanisation $F_{\text{Urb}}$ is computed for every model grid square from the US-Navy data.

Figure 10.2  (a) Roughness length for vegetation $z_{0\text{mVeg}}$ as provided by Baumgartner et al. (1977) on a $5^\circ \times 5^\circ$ grid.

(b) The roughness length field including orographic effects (as used by the model) $z_{0\text{m}}$ at T511 resolution plotted on a $1^\circ \times 1^\circ$ grid.
The vegetation roughness length for momentum $z_{0v\text{eg}}$ is blended with the urbanization roughness length ($z_{0\text{mUrb}}$) using a blending height $h_U$. The blending-height concept is based on the idea that internal boundary layers merge at a certain height which depends on the horizontal scale of the inhomogeneities (Claussen 1990). Below the blending height the different areas have their own logarithmic profile and the resulting stresses are averaged proportional to the area covered by the different roughness lengths. An effective roughness is defined such that over a homogeneous surface it would provide the same drag as the average over the two different surfaces. The effective roughness length $z_{0v\text{eg}}$ due to fraction $F_{\text{Urb}}$ with roughness length $z_{0\text{mUrb}}$ and fraction $(1 - F_{\text{Urb}})$ with roughness length $z_{0\text{mVeg}}$ can be found by solving:

$$\frac{1}{\left[ \ln\left( \frac{h_{0\text{Veg}}}{z_{0\text{mVeg}} + 1} \right) \right]^2} = \frac{F_{\text{Urb}}}{\left[ \ln\left( \frac{h_{0\text{Veg}}}{z_{0\text{mUrb}} + 1} \right) \right]^2} + \frac{1 - F_{\text{Urb}}}{\left[ \ln\left( \frac{h_{0\text{Veg}}}{z_{0\text{mVeg}} + 1} \right) \right]^2}$$

The blended roughness length for heat $z_{0h\text{Veg}}$ is computed from

$$\ln\left( \frac{h_{0\text{Veg}}}{z_{0\text{mVeg}} + 1} \right) = \ln\left( \frac{h_{0\text{Veg}}}{z_{0\text{Veg}} + 1} \right) \frac{\ln\left( \frac{h_{0\text{Veg}}}{z_{0\text{mVeg}} + 1} \right)}{\ln\left( \frac{h_{0\text{Veg}}}{z_{0\text{Veg}} + 1} \right)},$$

where $h_{0\text{Urb}} = 100$ m, $z_{0h\text{Veg}} = z_{0\text{mVeg}}/10$ and $z_{0\text{mUrb}} = 2.5$ m.

To compute the orographic contribution to the roughness lengths, a slope parameter ($S_L$) is needed and a characteristic height ($\hat{h}$) of the subgrid orography. From these the typical horizontal scale of
the subgrid orography can be derived ($\lambda$). Because the horizontal scales up to 10 km are the most important ones for the roughness lengths, we use also US-Navy information about maximum height, minimum height and number of significant ridges inside the $30' \times 30'$ squares. Two contributions to the subgrid standard deviation are computed (i.e. subgrid to the ECMWF model): the standard deviation $\sigma_i$ resolved by the $30' \times 30'$ data and the standard deviation $\sigma_s$ subgrid to the $30' \times 30'$ data

$$
\sigma_i^2 = \sum_i p_i h_i^2 - \left( \sum_i p_i h_i \right)^2
$$

$$
\sigma_s^2 = \sum_i \frac{1}{4} (h_i - h_{i\text{min}})(h_{i\text{max}} - h_i)
$$

$$
\sigma^2 = \sigma_i^2 + \sigma_s^2
$$

where:

$N$ = Number of relative $h_i$ maxima in the ECMWF grid square,

$F$ = Surface area of the ECMWF grid square,

$n_i$ = Number of significant ridges in the $i$th 10′ grid square,

$h_i$ = Mean height in the $i$th 10′ grid square,

$h_{i\text{max}}$ = Maximum height in the $i$th 10′ grid square,

$h_{i\text{min}}$ = Minimum height in the $i$th 10′ grid square,

$f_i$ = Surface area of the $i$th 10′ grid square,

$p_i$ = Proportion of the ECMWF grid square occupied by the $i$th 10′ grid square.

Mason (1991) uses slope parameter $S_i = \sum A/F$, where the summation is over all obstacles in area $F$ (an ECMWF grid square) and $A$ is the frontal, or wind-swept area, of the obstacles. The swept area of the resolved part is estimated from $\sigma_i$ and the number of relative maxima in the ECMWF grid square $N$. We assume that the height difference between valleys and hill tops is about $4\sigma_i$. So the resolved part of $\sum A/F$ is (assuming that the individual hills have a vertical dimension of $4\sigma_i$ and a horizontal dimension of $\sqrt{F/N}$ and that there are $N$ hills in a grid square):

$$
\left( \frac{\sum A}{F} \right)_i = \frac{4\sigma_i \sqrt{F/N}}{F}
$$

$$
= 4\sigma_i \sqrt{N/F}
$$

Similarly for the unresolved part:
\[
\left( \frac{\sum A}{F} \right)_i = \frac{\sum 4\sigma_i \sqrt{f_i / n_i n_i}}{F} \\
= \sum_i 4\sigma_i \sqrt{\frac{n_i f_i}{F}} \\
= \sum_i 4\sigma_i \sqrt{\frac{n_i}{f_i}} p_i
\]

where

\[p_i = \frac{f_i}{F}\]
\[\sigma_i^2 = \frac{1}{4} (h_i - h_i^{\text{min}})(h_i^{\text{max}} - h_i)\]

The total \(\sum A/F\) is:
\[
S_f = \frac{\sum A}{F} = \left( \frac{\sum A}{F} \right)_f + \left( \frac{\sum A}{F} \right)_s .
\]

(c) For the computation of the effective roughness lengths, two different formulations are used in principle: the gentle orography approximation (Taylor 1987) and the steep orography formulation (Mason 1991). For the gentle orography approximation we need the horizontal wavelength \(\lambda\) of the subgrid terrain which we estimate as:

\[\lambda = \frac{h}{S_f}, \text{ where } h = 4\sigma .\]

The effective roughness lengths \(z_{0m}\) and \(z_{0h}\) for gentle slopes is determined by:

\[\ln\left( \frac{\lambda}{2\pi z_{0m}} + 1 \right) = \ln\left( \frac{\lambda}{2\pi z_{0m\text{Veg1}}} + 1 \right) + C_e \left( \frac{S_f}{\pi} \right)^{\frac{1}{2}}\]
\[\ln\left( \frac{\lambda}{2\pi z_{0h}} + 1 \right) = \ln\left( \frac{\lambda}{2\pi z_{0h\text{Veg1}}} + 1 \right) + \frac{\ln\left( \frac{\lambda}{2\pi z_{0m\text{Veg1}}} + 1 \right)}{\ln\left( \frac{\lambda}{2\pi z_{0m}} + 1 \right)}\]

with \(C_e = 5\). For steep slopes the effective roughness lengths \(z_{0m}\) and \(z_{0h}\) are computed from:
\[
\ln \left( \frac{h}{2\sigma_{0m}} + 1 \right) = \left( \frac{k^2}{\frac{1}{2}C_d S_f + \left( \frac{k}{\ln(h/2\sigma_{0mVeeg1}) + 1} \right)^2} \right)^{\frac{1}{2}} \\
\ln \left( \frac{h}{2\sigma_{0h}} + 1 \right) = \ln \left( \frac{h}{2\sigma_{0hVeeg1}} + 1 \right) \frac{\ln \left( \frac{h}{2\sigma_{0mVeeg1}} + 1 \right)}{\ln \left( \frac{h}{2\sigma_{0m}} + 1 \right)}
\]

where \( C_d = 0.4 \), \( k = 0.4 \) and the transition from gentle to steep formulation should be at about \( S_f = 0.2/\pi \). In practise the quality of the US-Navy data is not sufficient to resolve the gentle slopes, so the treshold is set to 0 which implies that the steep formulation is always used. Another disadvantage of the transition from the gentle to steep slope formulation is that it is discontinuous, which reflects the sudden transition from attached to separated flow.

Orographic corrections are not applied for \( h < 10 \text{ m} \) and for \( S_f < 0.01 \). \( z_{0m} \) is not allowed to become larger than \( h/10 \) and is clipped at 100 m; \( z_{0h} \) has an imposed lower bound of \( e^{-20} \).

\[(d)\] A Gaussian filter (the same as for the mean orography) is applied to \( \ln(z_{0m}) \) and \( \ln(z_{0h}) \), the sea points are reset to \( z_{0m} = 0.001 \) and \( z_{0h} = 0.0001 \). Because of the wide dynamical range of parameter \( z_{0h} \), \( \ln(z_{0h}) \) is GRIB-coded and provided as input to the model.

### 10.6 Parameters for Gravity-Wave and Orographic Drag Schemes

The following subgrid parameters are needed: standard deviation \( \mu_{GW} \), anisotropy \( \gamma_{GW} \), orientation \( \theta_{GW} \), and slope \( \sigma_{GW} \). They are computed as follows (see [Lott and Miller 1997; Baines and Palmer 1990]):

\[(a)\] For every point (index \( i \)) of the 2°30’ data, \( \partial h/\partial x \) and \( \partial h/\partial y \) are computed by central differencing with help of the points to the north, south, east and west. These derivatives are computed after subtracting the mean orography at target resolution to avoid contributions from the slope of the resolved orography. The central differences in the North South direction use adjacent points; derivatives in the East West direction use adjacent points in the tropics but use equidistant points rather than equi-latitude points when approaching the polar regions (to maintain a uniform resolution over the globe). Then parameters \( K \), \( L \), and \( M \) are computed by summation, taking into account the weights \( p_i \) of every \( 10' \times 10' \) area in the ECMWF grid:

\[
K = \frac{1}{2} \sum_i p_i \left[ \left( \frac{\partial h}{\partial x} \right)_i^2 + \left( \frac{\partial h}{\partial y} \right)_i^2 \right] \\
L = \frac{1}{2} \sum_i p_i \left[ \left( \frac{\partial h}{\partial x} \right)_i^2 - \left( \frac{\partial h}{\partial y} \right)_i^2 \right] \\
M = \sum_i p_i \left( \frac{\partial h}{\partial x} \right)_i \left( \frac{\partial h}{\partial y} \right)_i
\]

\[(b)\] Anisotropy \( \gamma_{GW} \), orientation \( \theta_{GW} \), and slope \( \sigma_{GW} \) are computed from \( K \), \( M \) and \( L \):
and the standard deviation $\mu_{GW}$:

$$\mu_{GW} = \sum_i p_i h_i^2 - \left( \sum_i p_i h_i \right)^2 .$$

No further filtering is applied to the fields. Results are shown in Figs. 10.4, 10.5, 10.6, and 10.7.

Figure 10.4 Anisotropy $\gamma_{GW}$ of subgrid orography (1 indicates isotropic, 0 means maximum anisotropy)
Figure 10.5 Orientation $\theta_{GW}$ of subgrid orography.

Figure 10.6 Slope $\sigma_{GW}$ of subgrid orography.
Vegetation is represented by 4 climatological parameters: vegetation cover of low vegetation, vegetation cover of high vegetation, low vegetation type and high vegetation type. These parameters are derived from the 2’30” GLCC data by averaging over the target grid squares. The fractional covers for low and high vegetation are obtained by combining the fractions from all the low and high vegetation types of Table 10.1. The index of the dominant low and high vegetation types are also coded as climatological fields for use by the land surface scheme. The latter two fields can not be interpolated by standard procedures to another resolution. The resulting fields are shown in Fig. 10.8, Fig. 10.9, Fig. 10.10 and Fig. 10.11. Table 10.2 and Table 10.3 contain statistical information on the number of points in each vegetation class.

**Table 10.2**  **Percentage of land points at T511 for each low vegetation type**

<table>
<thead>
<tr>
<th>Index</th>
<th>Vegetation type</th>
<th>Percentage of land points</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>Crops, Mixed Farming</td>
<td>22.1</td>
</tr>
<tr>
<td>2</td>
<td>Short Grass</td>
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<td>7</td>
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<td>13.0</td>
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<td>Tundra</td>
<td>8.9</td>
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<tr>
<td>10</td>
<td>Irrigated Crops</td>
<td>4.7</td>
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<tr>
<td>11</td>
<td>Semidesert</td>
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<tr>
<td>16</td>
<td>Evergreen Shrubs</td>
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Table 10.2  Percentage of land points at T511 for each low vegetation type

<table>
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<th>Vegetation type</th>
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<td></td>
<td>Remaining land points without low vegetation</td>
<td>19.7</td>
</tr>
</tbody>
</table>

Table 10.3  Percentage of land points at T511 for each high vegetation type

<table>
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<th>Vegetation type</th>
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<td>Evergreen Broadleaf Trees</td>
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<td>19</td>
<td>Interrupted Forest</td>
<td>29.6</td>
</tr>
<tr>
<td></td>
<td>Remaining land points without high vegetation</td>
<td>37.0</td>
</tr>
</tbody>
</table>

Figure 10.8  Fractional cover of low vegetation.
Figure 10.9 Fractional cover of high vegetation.

Figure 10.10 Low vegetation type
10.8 ALBEDO

The background land albedo is interpolated to the model grid from the monthly mean values of a snow-free albedo produced for the combined 1982-1990 years. The albedo for that dataset was computed using the method of Sellers et al. (1996), but with new maps of soil reflectance, new values of vegetation reflectance, and the biophysical parameters described by Los et al. (2000).

The fields for January, April, July and October are shown in Fig. 10.12, Fig. 10.13, Fig. 10.14, and Fig. 10.15. To obtain a smooth evolution in time, the model does a linear interpolation between successive months, assuming that the monthly field applies to the 15th of the month. The model adapts the background albedo over water, ice and snow as documented in the chapter on radiation.
Figure 10.12  Climatological background albedo for January.

Figure 10.13  Climatological background albedo for April.
10.9 AEROSOLS

Aerosols are considered in the model following Tanre et al. (1984). The continental, maritime, urban and desert aerosols are geographically distributed over ice-free land, open sea, industrialized area and desert using a bi-Gaus-
sian horizontal filter of radius 2000 km to get overlapping distributions of each aerosol type, with a maximum optical thickness of 0.2, 0.05, 0.1, and 1.9, respectively (see Figs. 10.16 (a)–(d) for the geographical distributions and Fig. 10.16 (e) for the corresponding profiles). Well-mixed (vertically and horizontally) tropospheric background aerosols with an optical thickness of 0.03 and stratospheric background aerosols with an optical thickness of 0.045 are added to the previous amounts with a rate of change of optical thickness with pressure of 0.037 and 0.233 /atm respectively. The transition from troposphere to stratosphere is obtained by multiplication of the background values with $1 - L_{\text{stratos}}$ and $L_{\text{stratos}}$ respectively (see Fig. 10.16 (e)).

Figure 10.16 Distribution of (a) maritime and (b) continental type aerosols.
Figure 10.16 Distribution of (c) urban and (d) desert type aerosols.
Figure 10.16 (e) Type 1 (full line) profiles apply to maritime, continental and urban type aerosols; type 2 (short dashed line) applies to desert type; the third curve (long dashed line) represents $L_{\text{stratos}}$ and is used to determine the transition from tropospheric to stratospheric background aerosols.

10.10 OZONE

The ozone climatology that is operational since August 1997, distributes the ozone mixing ratio as a function of pressure, latitude and month following Fortuin and Langematz (1995).

Figure 10.17 Ozone climatology prescribed as a zonal mean according to the climatology by Fortuin and Langematz (1994). Operational in the ECMWF model since August 1997 (mass mixing ratio $\times 10^6$ on 60 levels).
Part IV: PHYSICAL PROCESSES

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