# ITU-R <br> Radiocommunication Sector of ITU 

Report ITU-R P.2297-1<br>(05/2019)

# Electron density models and data for transionospheric radio propagation 

P Series
Radiowave propagation

## Foreword

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Note: This ITU-R Report was approved in English by the Study Group under the procedure detailed in Resolution ITU-R 1

## REPORT ITU-R P.2297-1 <br> Electron density models and data for transionospheric radio propagation

(2013-2019)

## Scope

This Report presents the detailed characteristics of electron density and total electron content models essential to Recommendation ITU-R P. 531 for transionospheric propagation and supplementary digital Vertical TEC grid point maps.

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## 1 Introduction

As reported in Recommendation ITU-R P.531, a number of transionospheric propagation effects, such as refraction, dispersion and group delay, are in magnitude directly proportional to the Total Electron Content (TEC); Faraday rotation is also approximately proportional to TEC, with the contributions from different parts of the ray path weighted by the longitudinal component of magnetic field. Knowledge of the TEC thus enables many important ionospheric effects to be estimated quantitatively.
For estimating TEC, a procedure based on NeQuick, is recommended in Recommendation ITU-R P.531. The NeQuick model is also suitable for slant TEC evaluation. NeQuick is a climatological electron density models and its details are provided in the following sections. The last section includes example Vertical TEC monthly mean (and standard deviation) grid maps data obtained from measurements at a set of periods and solar activities.

## 2 NeQuick2 electron density model

### 2.1 Background

NeQuick 2 is the latest version of the NeQuick ionosphere electron density model developed at the Aeronomy and Radiopropagation Laboratory of the Abdus Salam International Centre for Theoretical Physics (ICTP) - Trieste, Italy with the collaboration of the Institute for Geophysics, Astrophysics and Meteorology of the University of Graz, Austria. The NeQuick is a quick-run ionospheric electron density model particularly designed for transionospheric propagation applications.

To describe the electron density of the ionosphere up to the peak of the F2 layer, the NeQuick uses a profile formulation which includes five semi-Epstein layers with modelled thickness parameters. Three profile anchor points are used: the E layer peak, the F1 peak and the F2 peak, that are modelled in terms of the ionosonde parameters $f o E, f o F 1, f o F 2$ and $M(3000) F 2$. These values can be modelled (e.g. ITU-R coefficients for $f o F 2, M 3000$ ) or experimentally derived. A semi-Epstein layer represents the model topside with a height - dependent thickness parameter empirically determined. The NeQuick package includes routines to evaluate the electron density along any ray-path and the corresponding TEC by numerical integration.

### 2.2 The NeQuick electron density model

This section describes how to compute the NeQuick electron density for a given location (identified by the coordinates $\mathrm{h}, \varphi, \lambda$ ) at a given time (month, UT) using Rz or F10.7 parameters (provided or not by the user).

Inputs:
height $h(\mathrm{~km})$, latitude $\varphi$ (degrees), longitude $\lambda$ (degrees), month $m t h$, Universal Time $U T$ (hours and decimals), Rz or F10.7.

Output:
electron density $N\left(\mathrm{~m}^{-3}\right)$.

### 2.2.1 The Epstein function

Since to compute the NeQuick electron density profile, the Epstein function is used, its formulation is recalled as:

$$
\begin{equation*}
\operatorname{Epst}(X, Y, Z, W)=\frac{X \exp \left(\frac{W-Y}{Z}\right)}{\left(1+\exp \left(\frac{W-Y}{Z}\right)\right)^{2}} \tag{1}
\end{equation*}
$$

### 2.2.2 Constants used

For the calculation of the slant TEC, some constant parameters are used and they are summarized in Table 1.

TABLE 1

## Constants definition

| Symbol | Constant description | Value | Units |
| :---: | :---: | :---: | :---: |
| DR | Degree to radiant conversion factor | $\pi / 180$ | $\mathrm{rad} / \mathrm{degrees}$ |
| RD | Radiant to degree conversion factor | $180 / \pi$ | $\mathrm{rad} / \mathrm{degrees}$ |
| $\chi_{0}$ | Zenith angle at night-day transition | 86.23 | degrees |
| $\mathrm{R}_{\mathrm{E}}$ | Earth mean radius | 6371.2 | km |

### 2.2.3 Auxiliary parameters

To compute the NeQuick electron density, several auxiliary parameters are preliminarily evaluated through specific modules. In this section the parameter formulation is given.

### 2.2.3.1 Local Time

Compute local time $L T$ (in hours and decimals) for the location considered.
Inputs:
longitude $\lambda$ (degrees), Universal Time $U T$ (hours and decimals).
Output:
local time $L T$ (hours and decimals).

$$
\begin{equation*}
L T=U T+\lambda / 15 \tag{2}
\end{equation*}
$$

### 2.2.3.2 Modip map

NeQuick 2.1 uses a map of modified dip or Modip to get the magnetic coordinates needed by the model for computing the F2 layer peak and F1 layer height.
Modip, which was first introduced by Rawer [1963, RD4] is defined by:

$$
\begin{equation*}
\tan \mu=\frac{I}{\sqrt{\cos \varphi}} \tag{3}
\end{equation*}
$$

in which $I$ is magnetic inclination (degrees) at 300 km and $\varphi$ is the latitude (degrees) of the considered location. NeQuick 2.1 model uses a grid of modip values contained in the file modip.asc.
The file modip.asc contains the values of modip on a geocentric grid from $90^{\circ} \mathrm{S}$ to $90^{\circ} \mathrm{N}$ with steps of 1 degree in latitude and from $180^{\circ} \mathrm{W}$ to $180^{\circ} \mathrm{E}$ with steps of 2 degrees in longitude. In particular the first 181 numbers are the modip values for latitude $($ degrees $)=-90.0($ and longitude $($ degrees $)=-$
$180,-178 \ldots, 180$ ), etc.; the last 181 numbers are the modip values for latitude (degrees) $=+90.0$ (and longitude (degrees) $=-180,-178 \ldots, 180)$.

To compute the modip value in a point at latitude $\varphi$ and longitude $\lambda$, a third order interpolation is applied using surroundings grid values.
The NeQuick interpolation function is described in § 2.3.3.1.
For computational purposes, it is therefore convenient to describe how the modip values are stored in a grid $D$ and how additional rows and columns are defined in order to apply the interpolation algorithm near the poles or near longitude $($ degrees $)=180$.

### 2.2.3.2.1 Store the modip.asc values in an array

$$
\begin{equation*}
D: \quad d_{i, j} \quad i=1, \ldots, 181 ; j=0, \ldots, 180 \tag{4}
\end{equation*}
$$

being $i$ the index of latitude and $j$ the index of longitude.
Define additional rows:

$$
\begin{align*}
d_{0, j} & =d_{2, \bmod (j+90,180)}  \tag{5}\\
d_{182, j} & =d_{180, \bmod (j+90,180)}  \tag{6}\\
d_{183, j} & =d_{179, \bmod (j+90,180)} \tag{7}
\end{align*}
$$

with $j=0, \ldots, 180$ and being $\bmod (\mathrm{a}, \mathrm{b})$ the remainder on division of a by b .
Define additional columns:

$$
\begin{gather*}
d_{i,-1}=d_{i, 179}  \tag{8}\\
d_{i, 181}=d_{i, 1} \tag{9}
\end{gather*}
$$

with $i=0, \ldots, 183$.
Therefore now the matrix $D$ becomes:

$$
\begin{equation*}
D: \quad d_{i, j} \quad i=0, \ldots, 183 ; j=-1, \ldots, 181 ; \tag{10}
\end{equation*}
$$

### 2.2.3.2.2 Compute the Modip

Inputs:
latitude $\varphi$ (degrees), longitude $\lambda$ (degrees), array $D$ (of dip latitude values).
Output:
Modip $\mu$ (degrees)
The selection of the interpolation grid-points is done by computing:

$$
\begin{equation*}
l=\operatorname{int}\left(\frac{\lambda+180}{2}\right)-2 \tag{11}
\end{equation*}
$$

If $l<0$ use

$$
\begin{equation*}
l=l+180 \tag{12}
\end{equation*}
$$

If $l>177$ use

$$
\begin{equation*}
l=l-180 \tag{13}
\end{equation*}
$$

Compute

$$
\begin{align*}
& a=\frac{\varphi+90}{1}+1  \tag{14}\\
& x=a-\operatorname{int}(a)  \tag{15}\\
& i=\operatorname{int}(a)-2 \tag{16}
\end{align*}
$$

For $k=1,4$; for $j=1,4$ build the $z_{j, k}$ as:

$$
\begin{equation*}
z_{j, k}=d_{i+j, l+k} \tag{17}
\end{equation*}
$$

For $k=1,4$ compute

$$
\begin{equation*}
z_{k}=z_{x}\left(z_{1, k}, z_{2, k}, z_{3, k}, z_{4, k}, x\right) \tag{18}
\end{equation*}
$$

using the interpolation function described in § 2.3.3.1.
Finally compute

$$
\begin{gather*}
b=\frac{\lambda+180}{2}  \tag{19}\\
y=b-\operatorname{int}(b) \tag{20}
\end{gather*}
$$

and using the interpolation function described in $\S$ 2.3.3.1, calculate

$$
\begin{equation*}
\mu=z_{x}\left(z_{1}, z_{2}, z_{3}, z_{4}, y\right) \tag{21}
\end{equation*}
$$

### 2.2.3.3 Solar declination

Compute $\sin \delta \delta_{s u n}, \cos \delta s u n$, the sine and cosine of the solar declination.
Inputs:
month $m t h$, Universal Time $U T$ (h)
Outputs:
$\sin \delta_{s u n}, \cos \delta \delta_{s u n}$,
Compute day of year at the middle of the month:

$$
\begin{equation*}
d_{y}=30.5 m t h-15 \tag{22}
\end{equation*}
$$

Compute time (days):

$$
\begin{equation*}
t=d_{y}+(18-U T) / 24 \tag{23}
\end{equation*}
$$

Compute the argument:

$$
\begin{equation*}
a_{m}=(0.9856 t-3.289) D R \tag{24}
\end{equation*}
$$

$$
\begin{equation*}
a_{l}=a_{m}+\left[1.916 \sin \left(a_{m}\right)+0.020 \sin \left(2 a_{m}\right)+282.634\right] D R \tag{25}
\end{equation*}
$$

Finally compute sine and cosine of solar declination:

$$
\begin{align*}
& \sin \delta_{S u n}=0.39782 \sin a_{l}  \tag{26}\\
& \cos \delta_{S u n}=\sqrt{1-\sin ^{2} \delta_{S u n}} \tag{27}
\end{align*}
$$

### 2.2.3.4 Solar zenith angle

Compute solar zenith angle $\chi$ (degrees) for the given location.
Inputs:
latitude $\varphi$ (degrees), local time $L T(\mathrm{~h}), \sin \delta_{s u n}, \cos \delta \delta_{s u}$,
Output:
solar zenith angle $\chi$ (degrees).
Compute

$$
\begin{gather*}
\cos _{\chi}=\sin (\phi D R) \sin \delta_{S u n}+\cos (\phi D R) \cos \delta_{S u n} \cos \left(\frac{\pi}{12}(12-L T)\right)  \tag{28}\\
\chi=R D \operatorname{atan} 2\left(\sqrt{1-\cos \chi^{2}, \cos \chi}\right) \tag{29}
\end{gather*}
$$

### 2.2.3.5 Effective solar zenith angle

Compute the effective solar zenith angle $\chi_{\text {eff }}$ (degrees) as a function of the solar zenith angle $\chi$ (degrees) and the solar zenith angle at day night transition $\chi_{0}$ (degrees).
Inputs:
Solar zenith angle $\chi$ (degrees), $\chi_{0}$ (degrees)
Output:
Effective solar zenith angle $\chi_{\text {eff }}$ (degrees).
Being

$$
\begin{equation*}
\chi_{0}=86.23^{\circ} \tag{30}
\end{equation*}
$$

Then

$$
\begin{equation*}
\chi_{0}=\frac{\chi+[90-0.24 \exp (20-0.2 \chi)] \exp \left[12\left(\chi-\chi_{0}\right)\right]}{1+\exp \left[12\left(\chi-\chi_{0}\right)\right]} \tag{31}
\end{equation*}
$$

### 2.2.4 Model parameters

In the following point, model peak parameter and auxiliary parameter values will be calculated.

### 2.2.4.1 $\quad f o E$ and $N m E$

To compute the E layer critical frequency $f o E(\mathrm{MHz})$ at a given location, in addition to the effective solar zenith angle $\chi_{\text {eff }}$, a season dependent parameter has to be computed.

Inputs:
latitude $\varphi$ (degrees), flux $f l x$, effective solar zenith angle $\chi_{\text {eff }}$ (degrees), month $m t h$.
Output:
foE (MHz).
Define the seas parameter as a function of the month of the year as follows:

$$
\begin{align*}
& \text { If } m t h=1,2,11,12 \text { then seas }=-1  \tag{32}\\
& \text { If } m t h=3,4,9,10 \text { then seas }=0  \tag{33}\\
& \text { If } m t h=5,6,7,8 \text { then seas }=1 \tag{34}
\end{align*}
$$

Introduce the latitudinal dependence:

$$
\begin{gather*}
e e=\exp (0.3 \varphi)  \tag{35}\\
\text { seasp }=\text { seas } \frac{e e-1}{e e+1} \tag{36}
\end{gather*}
$$

$$
\begin{equation*}
f o E=\sqrt{(1.112-0.019 \text { seasp })^{2} \sqrt{f l x}\left[\cos \left(\chi_{\mathrm{eff}} D R\right)\right]^{0.6}+0.49} \tag{37}
\end{equation*}
$$

The E layer maximum density $\operatorname{NmE}\left(10^{11} \mathrm{~m}^{-3}\right)$ as a function of $f o E(\mathrm{MHz})$ is computed as:

$$
\begin{equation*}
N m E=0.124 f o E^{2} \tag{38}
\end{equation*}
$$

### 2.2.4.2 foF1 and NmF 1

The $F 1$ layer critical frequency $f o F 1(\mathrm{MHz})$ in NeQuick 2.1 has been reformulated [Leitinger et al., 2005, RD3].

Inputs:
E layer critical frequency $f o E(\mathrm{MHz}), F 2$ layer critical frequency $f o F 2(\mathrm{MHz})$, Output:

> foF1 (MHz).

$$
\begin{gather*}
0  \tag{39}\\
f o F 1=
\end{gather*}=\left\{\begin{array}{cc}
f o E<2 \\
1.4 f o E & \text { if } \\
0.85 \cdot 1.4 f o E & \text { if }
\end{array} \quad 1.4 f o E>2 \text { and } 1.4 \text { foE } \leq 0.85 f o F 2\right.
$$

The F layer maximum density $\operatorname{NmF1}\left(10^{11} \mathrm{~m}^{-3}\right)$ as a function of $f o F 1(\mathrm{MHz})$ is computed as:

$$
\begin{equation*}
N m F 1=0.124 * f o F 1^{2} \tag{40}
\end{equation*}
$$

### 2.2.4.3 foF2 and NmF2; M(3000)F2

To compute $f o F 2$ and $M(3000) F 2$, the NeQuick model uses the ITU-R (formerly CCIR) coefficients [ITU-R, 1997]. These coefficients are stored in the ccirXX.asc files and are the spherical harmonic coefficients representing the development of monthly median $f o F 2$ and $M(3000) F 2$ all over the world. The coefficients correspond to low $\left(R_{12}=0\right)$ and high $\left(R_{12}=100\right)$ solar activity conditions. Therefore they have to be interpolated (or extrapolated) to obtain the set of coefficients corresponding to the required solar activity.
Each file ccirXX.asc contains 2858 values sequentially organized as follows: $\left\{f 2_{1,1,1}, f 2_{1,1,2}, \ldots, f 2_{1,1,13}\right.$, $f 2_{1,2,1}, f 2_{1,2,2}, \ldots, f 2_{1,2,13}, \ldots, f 2_{1,76,1}, f 2_{1,76,2}, \ldots, f 2_{1,76,13}, f 2_{2,1,1}, \ldots, f 2_{2,1,2}, \ldots, f 2_{2,1,13}, f 2_{2,2,1}, f 2_{2,2,2}, \ldots$, $f 2_{2,2,13}, \ldots, f 2_{2,76,1}, f 2_{2,76,2}, \ldots, f 2_{2,76,13}, f m 3_{1,1,1}, f m 3_{1,1,2}, \ldots, f m 3_{1,1,9}, f m 3_{1,2,1}, f m 3_{1,2,2}, \ldots, f m 3_{1,2,9}, \ldots$, $f m 3_{1,49,1}, f m 3_{1,49,2}, \ldots, f m 3_{1,49,9}, f m 3_{2,1,1}, f m 3_{2,1,2}, \ldots, f m 3_{2,1,9}, f m 3_{2,2,1}, f m 3_{2,2,2}, \ldots, f m 3_{2,2,9}, \ldots, f m 3_{2,49,1}$, $\left.f m 3_{2,49,2}, \ldots, f m 3_{2,49,9}\right\}$. (The notation used will become self evident after the definition of the F2 and Fm3 arrays).
The coefficients of a ccirXX. asc file have to be stored in two 3-D arrays, $F 2$ and $F m 3$, as indicated in the next section.

### 2.2.4.3.1 Store ccirXX. asc values

Input:
Month mth
Outputs:

## F2, Fm3

Select the file name to read:

$$
\begin{equation*}
X X=m+10 \tag{41}
\end{equation*}
$$

(e.g. ccir21.asc for November) and store the file content in the two arrays of coefficients:
coefficients for $f o F 2$

$$
\begin{equation*}
F 2: \quad f 2_{i, j, k} \quad i=1,2 ; j=1, \ldots, 76 ; k=1, \ldots, 13 \tag{42}
\end{equation*}
$$

coefficients for $M(3000) F 2$

$$
\begin{equation*}
\text { Fm3: } \quad f m 3_{i, j, k} \quad i=1,2 ; j=1, \ldots, 49 ; k=1, \ldots, 9 \tag{43}
\end{equation*}
$$

### 2.2.4.3.2 Interpolate ITU-R coefficients for $\boldsymbol{R}_{12}$

Compute $A F 2$, the array of interpolated coefficients for $f o F 2$ and $A m 3$, the array of interpolated coefficients for $M(3000) F 2$.
Inputs:
F2, Fm3, $R_{12}$
Outputs:
AF2, Am3
Compute the array of interpolated coefficients for $f o F 2$ :

$$
\begin{equation*}
A F 2: \quad \text { af } 2_{j, k} \quad j=1, \ldots, 76 ; k=1, \ldots, 13 \tag{44}
\end{equation*}
$$

$A F 2$ elements are calculated by linear combination of the elements of $F 2$ :

$$
\begin{equation*}
\text { af } 2_{j, k}=f 2_{1, j, k}\left(1-\frac{R_{12}}{100}\right)+f 2_{2, j, k} \frac{R_{12}}{100} \quad j=1, \ldots, 76 ; k=1, \ldots, 13 \tag{45}
\end{equation*}
$$

Compute the array of interpolated coefficients for $M(3000) F 2$ :

$$
\begin{equation*}
\text { Am3: } a m 3_{j, k} \quad j=1, \ldots, 49 ; k=1, \ldots, 9 \tag{46}
\end{equation*}
$$

Am3 elements are calculated by linear combination of the elements of Fm3:

$$
\begin{equation*}
a m 3_{j, k}=f m 3_{1, j, k}\left(1-\frac{R_{12}}{100}\right)+f m 3_{2, j, k} \frac{R_{12}}{100} \quad j=1, \ldots, 49 ; k=1, \ldots, 9 \tag{47}
\end{equation*}
$$

### 2.2.4.3.3 Compute Fourier time series for $f o F 2$ and $M(3000) F 2$

Inputs:
Universal Time $U T$ (h), arrays of interpolated ITU-R coefficients AF2, Am3
Outputs:
$C F 2, C m 3$, vectors of coefficients for Legendre calculation for $f o F 2$ and $M(3000) F 2$.
Vector CF2 has 76 elements:

$$
\begin{equation*}
C F 2: \quad c f 2_{l} \quad l=1, \ldots, 76 \tag{48}
\end{equation*}
$$

Vector Cm3 has 49 elements:

$$
\begin{equation*}
\text { Cm3: } \quad c m 3_{l} \quad l=1, \ldots, 49 \tag{49}
\end{equation*}
$$

Compute the time argument:

$$
\begin{equation*}
T=(15 U T-180) D R \tag{50}
\end{equation*}
$$

For $i=1, . ., 76$ calculate the Fourier time series for $f o F 2$ :

$$
\begin{equation*}
c f 2_{i}=a f 2_{i, 1}+\sum_{k=1}^{6}\left[a f 2_{i, 2 k} \sin (k T)+a f 2_{i, 2 k+1} \cos (k T)\right] \tag{51}
\end{equation*}
$$

For $i=1, . ., 49$ calculate the Fourier time series for $M(3000) F 2$ :

$$
\begin{equation*}
c m 3_{i}=a m 3_{i, 1}+\sum_{k=1}^{4}\left[a m 3_{i, 2 k} \sin (k T)+a m 3_{i, 2 k+1} \cos (k T)\right] \tag{52}
\end{equation*}
$$

### 2.2.4.3.4 Compute $f o F 2$ and $M(3000) F 2$ by Legendre calculation

Inputs:
Modip $\mu$ (degrees), latitude $\varphi$ (degrees), longitude $\lambda$ (degrees), vector $C F 2$ of the coefficients for Legendre combination for foF2, vector Cm3 of the coefficients for Legendre combination for $M(3000) F 2$.
Outputs:
foF2 (MHz), M(3000)F2
Define vectors containing sine and cosines of coordinates:

$$
\begin{equation*}
M: \quad m_{k} \quad k=1, \ldots, 12 \tag{53}
\end{equation*}
$$

$$
\begin{array}{lll}
P: & p_{n} & n=2, \ldots, 9 \\
& & \\
S: & s_{n} & n=2, \ldots, 9  \tag{56}\\
C: & c_{n} & n=2, \ldots, 9
\end{array}
$$

Compute modip coefficients:

$$
\begin{equation*}
m_{1}=1 \tag{57}
\end{equation*}
$$

and for $k=2, \ldots, 12$

$$
\begin{equation*}
m_{k}=\sin ^{k-1}(\mu D R) \tag{58}
\end{equation*}
$$

Compute latitude and longitude coefficients:
and for $n=2, \ldots, 9$

$$
\begin{gather*}
p_{n}=\cos ^{n-1}(\phi D R)  \tag{59}\\
s_{n}=\sin ((n-1) \lambda D R)  \tag{60}\\
c_{n}=\cos ((n-1) \lambda D R) \tag{61}
\end{gather*}
$$

Compute foF2
Order 0 term:

$$
\begin{equation*}
f o F 2_{1}=\sum_{k=1}^{12} c f 2_{k} m_{k} \tag{62}
\end{equation*}
$$

having the increased Legendre grades for $f o F 2$ in a vector:

$$
\begin{align*}
& Q: \quad q_{n} \quad n=1, \ldots, 9  \tag{63}\\
& Q=(12,12,9,5,2,1,1,1,1) \tag{64}
\end{align*}
$$

for computational need define also:

$$
\begin{align*}
K: & k_{n} \quad n=1, \ldots, 9  \tag{65}\\
k_{1}= & q_{1} \tag{66}
\end{align*}
$$

and for $n=2, \ldots, 9$

$$
\begin{equation*}
k_{n}=k_{n 1}+2 q_{n 1} \tag{67}
\end{equation*}
$$

for $n=2, \ldots, 9$ compute the higher order terms:

$$
\begin{equation*}
f o F 2_{n}=\sum_{k=1}^{q_{n}}\left(c f 2_{k_{n}}+2 \mathrm{k}-1 c_{n}+c f 2_{k_{n}}+2 \mathrm{k} s_{n}\right) m_{k} p_{n} \tag{68}
\end{equation*}
$$

Finally sum the terms to obtain $f o F 2$ :

$$
\begin{equation*}
f o F 2=\sum_{n=1}^{9} f o F 2_{n} \tag{69}
\end{equation*}
$$

Compute $M(3000) F 2$
Order 0 term:

$$
\begin{equation*}
M(3000) F 2_{0}=\sum_{k=1}^{5} c m 3_{k} m_{k} \tag{70}
\end{equation*}
$$

having the increased Legendre grades for $M(3000) F 2$ in a vector:

$$
\begin{gather*}
R: \quad r_{n} \quad n=1, \ldots, 7  \tag{71}\\
R=(7,8,6,3,2,1,1) \tag{72}
\end{gather*}
$$

for computational need define also:

$$
\begin{array}{r}
H: \quad h_{n} \quad n=1, \ldots, 7 \\
h_{1}=q_{1} \tag{74}
\end{array}
$$

and for $n=2, \ldots, 7$

$$
\begin{equation*}
h_{n}=h_{n 1}+2 q_{n 1} \tag{75}
\end{equation*}
$$

for $n=2, \ldots 7$, compute the higher order terms:

$$
\begin{equation*}
M(3000) F 2_{n}=\sum_{k=1}^{r_{n}}\left(c m 3_{h_{n}+2 k-1} c_{n}+c m 3_{h_{n}+2 k} s_{n}\right) m_{k} p_{n} \tag{76}
\end{equation*}
$$

Finally sum the terms:

$$
\begin{equation*}
M(3000)=\sum_{n=1}^{7} M(3000) F 2_{n} \tag{77}
\end{equation*}
$$

To compute NmF 2 use:

$$
\begin{equation*}
N m F 2=0.124 f o F 2^{2} \tag{78}
\end{equation*}
$$

where $N m F 2$ is in $\left(10^{11} \mathrm{~m}^{-3}\right)$.

### 2.2.4.4 hmE

The E layer maximum density height $h m E(\mathrm{~km})$ is defined as a constant:

$$
\begin{equation*}
h m E=120 \tag{79}
\end{equation*}
$$

### 2.2.4.5 hmF2

Compute the $F 2$ layer maximum density height $h m F 2(\mathrm{~km})$.
Inputs:
$f o E(\mathrm{MHz}), f o F 2(\mathrm{MHz}), M(3000) F 2$.
Output:
$h m F 2(\mathrm{~km})$.

$$
\begin{equation*}
h m F 2=\frac{1490 M \sqrt{\frac{0.0196 M^{2}+1}{1.2967 M^{2} 1}}}{M+M} 176 \tag{80}
\end{equation*}
$$

where:

$$
\begin{gather*}
M=M(3000) F 2  \tag{81}\\
\Delta M=-0.012 \quad \text { if } \quad f o E<10^{-30}  \tag{82}\\
\Delta M=\frac{0.253}{\rho-1.215}-0.012 \quad \text { if } \quad \text { fo } E \geq 10^{-30} \tag{83}
\end{gather*}
$$

and the ratio $\rho$ is computed as:

$$
\begin{equation*}
\rho=\frac{\frac{f o F 2}{f o E} \exp \left[20\left(\frac{f o F 2}{f o E}-1.75\right)\right]+1.75}{\exp \left[20\left(\frac{f o F 2}{f o E}-1.75\right)\right]+1} \tag{84}
\end{equation*}
$$

### 2.2.4.6 hmF1

Compute the $F 1$ layer maximum density height $h m F 1(\mathrm{~km})$ :
Inputs:
$h m E(\mathrm{~km}), h m F 2(\mathrm{~km})$.
Output:
$h m F 1(\mathrm{~km})$.

$$
\begin{equation*}
h m F 1=\frac{h m E+h m F 2}{2} \tag{85}
\end{equation*}
$$

### 2.2.4.7 $\quad$ B2 $2_{b o t}$

Compute the thickness parameter $B 2_{\text {bot }}(\mathrm{km})$.
Inputs:
$N m F 2\left(10^{11} \mathrm{~m}^{-3}\right)$, foF2 (MHz), $M(3000) F 2$.
Output:
$B 2_{\text {bot }}(\mathrm{km})$.

$$
\begin{equation*}
B 2_{\text {bot }}=\frac{0.385 \mathrm{NmF} 2}{0.01 \exp [-3.467+1.714 \ln (f o F 2)+2.02 \ln (M)]} \tag{86}
\end{equation*}
$$

where
$M=M(3000) F 2$ as already indicated in equation (82).

### 2.2.4.8 A1

Compute the $F 2$ layer amplitude $A 1\left(10^{11} \mathrm{~m}^{-3}\right)$.
Inputs:
$N m F 2\left(10^{11} \mathrm{~m}^{-3}\right)$.
Output:
A1 $\left(10^{11} \mathrm{~m}^{-3}\right)$.

$$
\begin{equation*}
A 1=4 N m F 2 \tag{87}
\end{equation*}
$$

### 2.2.4.9 $\quad$ A2

Compute the $F 1$ layer amplitude $A 2\left(10^{11} \mathrm{~m}^{-3}\right)$.
Inputs:
$N m F 1\left(10^{11} \mathrm{~m}^{-3}\right), A 1\left(10^{11} \mathrm{~m}^{-3}\right), h m F 2(\mathrm{~km}), B 2_{\text {bot }}(\mathrm{km}), h m F 1(\mathrm{~km})$.
Output:
A2 $\left(10^{11} \mathrm{~m}^{-3}\right)$.
First compute the auxiliary parameter

$$
\begin{equation*}
A 2 a=4.0\left(N m F 1-\operatorname{Epst}\left(A 1, h m F 2, B 2_{b o t}, h m F 1\right)\right) \tag{88}
\end{equation*}
$$

where the function Epst is the one defined in § 2.2.1.
Then compute

$$
\begin{equation*}
A 2=\frac{A 2 a \exp [60(A 2 a-0.005)]+0.05}{1+\exp [60(A 2 a-0.005)]} \tag{89}
\end{equation*}
$$

### 2.2.4.10 B1 $1_{\text {top }}$

Compute the thickness parameter $B 1_{\text {top }}(\mathrm{km})$.
Inputs:
$h m E(\mathrm{~km}), h m F 2(\mathrm{~km})$
Output:
$B 1_{\text {top }}(\mathrm{km})$.
The thickness parameter $B 1_{\text {top }}(\mathrm{km})$ is given by:

$$
\begin{equation*}
B 1_{\text {top }}=0.3(h m F 2-h m F 1) \tag{90}
\end{equation*}
$$

### 2.2.4.11 B1 $1_{\text {bot }}$

Compute the thickness parameter $B 1_{\text {bot }}(\mathrm{km})$.
Inputs:
$h m F 1(\mathrm{~km}), h m E(\mathrm{~km})$
Output:
$B 1_{\text {bot }}(\mathrm{km})$.

$$
\begin{equation*}
B 1_{b o t}=0.5(h m F 1-h m E) \tag{91}
\end{equation*}
$$

### 2.2.4.12 A3

Compute the E layer amplitude $A 3\left(10^{11} \mathrm{~m}^{-3}\right)$.
Inputs:
$\operatorname{NmE}\left(10^{11} \mathrm{~m}^{-3}\right), A 2\left(10^{11} \mathrm{~m}^{-3}\right), h m F 1(\mathrm{~km}), h m F 2(\mathrm{~km}), h m E(\mathrm{~km}), B 1_{b o t}(\mathrm{~km}), B 2_{\text {bot }}(\mathrm{km})$.
Output:
A3 ( $10^{11} \mathrm{~m}^{-3}$ ).
First compute the auxiliary parameter $A 3 a$ :

$$
\begin{equation*}
A 3 a=4.0\left(N m E-E p s t\left(A 2, h m F 1, B 1_{b o t}, h m E\right)-E p s t\left(A 1, h m F 2, B 2_{b o t}, h m E\right)\right) \tag{92}
\end{equation*}
$$

where the function Epst is the one defined in § 2.2.1.
Then compute:

$$
\begin{equation*}
A 3=\frac{A 3 a \exp [60(\mathrm{~A} 3 a-0.005)]+0.005}{1+\exp [60(\mathrm{~A} 3 a-0.005)]} \tag{93}
\end{equation*}
$$

### 2.2.4.13 $B E_{\text {top }}$

Compute the thickness parameter $B E_{\text {top }}(\mathrm{km})$.
Inputs:
$h m F 1(\mathrm{~km}), h m E$
Output:
$B E_{\text {top }}(\mathrm{km})$.

$$
\begin{equation*}
B E_{t o p}=\max (0.5(h m F 1-h m E), 7) \tag{94}
\end{equation*}
$$

### 2.2.4.14 $B E_{b o t}$

Compute the thickness parameter $B E_{\text {bot }}(\mathrm{km})$.

$$
\begin{equation*}
B E_{b o t}=5 \tag{95}
\end{equation*}
$$

### 2.2.4.15 Shape parameter $\boldsymbol{k}$

Compute the shape parameter $k$.
Inputs:
$f o F 2(\mathrm{MHz}), h m F 2(\mathrm{~km}), B 2_{\text {bot }}(\mathrm{km})$.
Output:
k.

Compute

$$
\begin{equation*}
k=3.22-0.0538 \text { foF } 2-0.00664 h m F 2+0.113 \frac{h m F 2}{B 2_{\text {bot }}}+0.00257 R_{12} \tag{96}
\end{equation*}
$$

a lower limiting value of 1 is imposed on $k$.

### 2.2.4.16 $\mathrm{H}_{0}$

Compute the topside thickness parameter $H_{0}(\mathrm{~km})$.
Inputs:
$B 2_{\text {bot }}(\mathrm{km}), k$
Output:
$H_{0}(\mathrm{~km})$
Compute

$$
\begin{equation*}
H_{0}=k B 2_{b o t} \tag{97}
\end{equation*}
$$

### 2.3 Electron density computation

To compute the electron density $N=N(h, \varphi, \lambda, l, f l x, m t h, U T)$ at a given point (identified by the coordinates $h, \varphi, \lambda$ ) and time (month, UT) using a given F10.7, all NeQuick parameters have to be evaluated. Nevertheless two different modules have to be used accordingly to the height considered. In particular
if

$$
\begin{equation*}
h \leq h m F 2 \tag{98}
\end{equation*}
$$

the bottomside electron density has to be computed using the algorithm illustrated in § 2.3.1.
if

$$
\begin{equation*}
h>h m F 2 \tag{99}
\end{equation*}
$$

the topside electron density has to be computed using the algorithm illustrated in § 2.3.2.

### 2.3.1 The bottomside electron density

Compute the electron density $N$ of the bottomside ( $h \leq h m F 2$ ).
Inputs:
height $h(\mathrm{~km}), A_{1}\left(10^{11} \mathrm{~m}^{-3}\right), A_{2}\left(10^{11} \mathrm{~m}^{-3}\right), A_{3}\left(10^{11} \mathrm{~m}^{-3}\right), h m F 2(\mathrm{~km}), h m F 1(\mathrm{~km}), h m E(\mathrm{~km}), B 2_{\text {bot }}$ $(\mathrm{km}), B 1_{\text {top }}(\mathrm{km}), B 1_{\text {bot }}(\mathrm{km}), B E_{\text {top }}(\mathrm{km}), B E_{\text {bot }}(\mathrm{km})$.

Output:
(bottomside) electron density $N\left(\mathrm{~m}^{-3}\right)$.
Select the relevant $B$ parameters for the current height:

$$
\begin{gather*}
B E= \begin{cases}B E_{\text {top }} & \text { if } h>h m E \\
B E_{\text {bot }} & \text { if } h \leq h m E\end{cases}  \tag{100}\\
B F 1= \begin{cases}B F 1_{\text {top }} & \text { if } h>h m F 1 \\
B F 1_{\text {bot }} & \text { if } h \leq h m F 1\end{cases} \tag{101}
\end{gather*}
$$

Compute the exponential arguments for each layer:

$$
\begin{gather*}
\alpha_{1}=\frac{h-h m F 2}{B 2_{b o t}}  \tag{102}\\
\alpha_{2}=\frac{h-h m F 1}{B F 1} \exp \left(\frac{10}{1+1|h-h m F 2|}\right)  \tag{103}\\
\alpha_{3}=\frac{h-h m E}{B E} \exp \left(\frac{10}{1+1|h-h m F 2|}\right) \tag{104}
\end{gather*}
$$

If $h<90 \mathrm{~km}$, for each $i=1,3$ compute:

$$
s_{i}= \begin{cases}\alpha_{i}=\alpha_{i} *(5+90-h) / 5 \\ 0 & \text { if }\left|\alpha_{i}\right|>25  \tag{106}\\ A_{i} \frac{\exp \left(\alpha_{i}\right)}{\left(1+\exp \left(\alpha_{i}\right)\right)^{2}} & \text { if }\left|\alpha_{i}\right| \leq 25\end{cases}
$$

Then compute the electron density as:

$$
\begin{equation*}
N=\left(s_{1}+s_{2}+s_{3}\right) \times 10^{11} \tag{107}
\end{equation*}
$$

### 2.3.2 The topside electron density

Compute the electron density $N$ of the topside ( $h>h m F 2$ ).
Inputs:
height $h(\mathrm{~km}), N m F 2\left(10^{11} \mathrm{~m}^{-3}\right), h m F 2(\mathrm{~km}), H_{0}(\mathrm{~km})$.
Output:
(topside) electron density $N\left(\mathrm{~m}^{-3}\right)$.
Define the constant parameters $g$ and $r$ as:

$$
\begin{gather*}
g=0.125  \tag{108}\\
r=100 \tag{109}
\end{gather*}
$$

compute the $z$ argument as:

$$
\begin{equation*}
z=\frac{h-h m F 2}{H_{0}\left[1+\frac{r g(h-h m F 2)}{r H_{0}+g(h-h m F 2)}\right]} \tag{110}
\end{equation*}
$$

Eventually

$$
\begin{equation*}
N(h)=\frac{4 N m F 2}{(1+\exp (z))^{2}} \exp (z) \tag{111}
\end{equation*}
$$

### 2.3.3 Auxiliary routines

### 2.3.3.1 Third order interpolation function $z_{x}\left(z_{1}, z_{2}, z_{3}, z 4, x\right)$

Be $P 1=\left(-1, z_{1}\right), P 2=\left(0, z_{2}\right), P 3=\left(1, z_{3}\right), P 4=\left(2, z_{4}\right)$ :
If $P=\left(x, z_{x}\right)$, to compute the interpolated value $z x$ at the position $x$, being $x \in[0,1)$, the following algorithm is applied.
Inputs: $z_{1}, z_{2}, z_{3}, z_{4}, x$
Outputs: $z x$.
If $\left|x^{2}\right| \leq 10^{10}$

$$
\begin{equation*}
z_{x}=z_{2} \tag{112}
\end{equation*}
$$

Otherwise compute:

$$
\begin{gather*}
\quad=2 x \quad 1  \tag{113}\\
g_{1}=z_{3}+z_{2}  \tag{114}\\
g_{2}=z_{3}-z_{2}  \tag{115}\\
g_{3}=z_{4}+z_{1}  \tag{116}\\
g_{4}=\frac{z_{4}-z_{1}}{3}  \tag{117}\\
a_{0}=9 g_{1}-g_{3}  \tag{118}\\
a_{1}=9 g_{2}-g_{4}  \tag{119}\\
a_{2}=g_{3}-g_{1}  \tag{120}\\
a_{3}=g_{4}-g_{2}  \tag{121}\\
z_{x}=\frac{1}{16}\left(a_{0}+a_{1} \delta+a_{2} \delta^{2}+a_{3} \delta^{3}\right) \tag{122}
\end{gather*}
$$

### 2.4 TEC calculation

To compute the slant TEC along a straight line between a point $P_{1}$ and a point $P_{2}$, the NeQuick electron density $N$ has to be evaluated on a point $P$ defined by the coordinates $\{h, \varphi, \lambda\}$ along the raypath.
It is a choice depending on receiver computation capabilities to identify the number of points where $N$ is to be evaluated, in order to obtain a sufficient accuracy for a subsequent integration, leading to slant TEC.

The Earth is assumed to be a sphere with a radius of 6371.2 km .
For computational efficiency, if the latitude of $P_{1}$ and $P_{2}$ are close to each other and the longitude of $P_{1}$ and $P_{2}$ are close to each other (as indicated in equation (127)):

$$
\left\{\begin{array}{l}
\left|\varphi_{2}-\varphi_{1}\right|<10^{-5}  \tag{123}\\
\left|\lambda_{2}-\lambda_{1}\right|<10^{-5}
\end{array}\right.
$$

the vertical integration algorithm has to be used, as described in § 2.4.1; otherwise use the slant integration algorithm as described in § 2.4.2.
During the TEC computation, the electron density at the point $P$ has to be evaluated as indicated in $\S 2.2$, while the calculation of the coordinates of the point $P$ along the ray-path will be described in $\S$ 2.4.1.1 if a vertical ray-path is considered and in § 2.4.2.3 if a slant ray-path is considered.

### 2.4.1 Vertical TEC calculation

To compute NeQuick vertical TEC, first compute all profile parameters $h m E, h m F 1, h m F 2, A 1, A 2$, $A 3, B 2_{\text {bot }}, B 1_{\text {top }}, B 1_{\text {bot }}, B E_{\text {top }}, B E_{\text {bot }}, N m F 2, H_{0}$, then compute the integration of the electron density (bottomside or topside) as a function of height:

$$
\begin{equation*}
T E C=\int N(h) d h \tag{124}
\end{equation*}
$$

being:

$$
\begin{align*}
& h_{1}=r_{1}-R_{E}  \tag{125}\\
& h_{2}=r_{2}-R_{E} \tag{126}
\end{align*}
$$

### 2.4.1.1 Vertical TEC numerical integration

Inputs:

- Integration endpoints $h_{1}(\mathrm{~km}), h_{2}(\mathrm{~km})$
- Integration accuracy $\varepsilon$
- Model parameters $A_{1}\left(10^{11} \mathrm{~m}^{-3}\right), A_{2}\left(10^{11} \mathrm{~m}^{-3}\right), A_{3}\left(10^{11} \mathrm{~m}^{-3}\right), h m F 2(\mathrm{~km}), h m F 1(\mathrm{~km}), h m E$ $(\mathrm{km}), B 2_{\text {bot }}(\mathrm{km}), B 1_{\text {top }}(\mathrm{km}), B 1_{\text {bot }}(\mathrm{km}), B E_{\text {top }}(\mathrm{km}), B E_{\text {bot }}(\mathrm{km}), N m F 2\left(\mathrm{~m}^{-3}\right), H_{0}(\mathrm{~km})$
Output:
- $\quad \operatorname{TEC}\left(10^{15} \mathrm{~m}^{-2}\right)$

Being $\varphi, \lambda$, and all model parameters fixed during the integration, in the following a simplified notation is used:

$$
N(h)= \begin{cases}\text { bottomside } N & \text { if } h \leq h m F 2  \tag{127}\\ \text { topside } N & \text { if } h>h m F 2\end{cases}
$$

$N(h)$ is computed using the algorithms described in § 2.3.
Start the calculation using 8 points:

$$
\begin{equation*}
n=8 \tag{128}
\end{equation*}
$$

Repeat the following computations until the desired accuracy is obtained.
Calculate the integration intervals:

$$
\begin{gather*}
\Delta_{n}=\frac{h_{2}-h_{1}}{n}  \tag{129}\\
g=0.5773502691896 \cdot \Delta_{n}  \tag{130}\\
y=g_{1}+\frac{\Delta_{n}-g}{2}  \tag{131}\\
G N_{2}=\frac{\Delta_{n}}{2} \sum_{i=0}^{n-1}\left[N\left(y+i \Delta_{n}\right)+N\left(y+i \Delta_{n}+g\right)\right] \tag{132}
\end{gather*}
$$

Double the number of points:

$$
\begin{equation*}
n=2 n \tag{133}
\end{equation*}
$$

and define

$$
\begin{equation*}
G N_{1}=G N_{2} \tag{134}
\end{equation*}
$$

repeating steps from (129) to (132) it is now possible to compare the two values obtained to see if the required accuracy is achieved: if

$$
\begin{equation*}
\left|G N_{1}-G N_{2}\right|>\varepsilon\left|G N_{1}\right| \tag{135}
\end{equation*}
$$

then continue increasing the number of points using (133), redefine $G N_{1}$ using (134) and repeat again steps from (129) to (132).

When the test (135) fails, the required accuracy has been reached, and the value of the integral is obtained by:

$$
\begin{equation*}
T E C=\left(G N_{2}+\frac{G N_{2}-G N_{1}}{15}\right) \times 10^{-13} \tag{136}
\end{equation*}
$$

### 2.4.2 Slant TEC calculation

To compute the electron density at a point $P$ along the slant ray-path defined by the points $P_{1}$ and $P_{2}$ the following specific geometrical configuration is considered.

### 2.4.2.1 Geometrical configuration

To simplify the formulation it is assumed that if $\alpha$ is an angle in (degrees), $\tilde{\alpha}$ is the same angle in (rad):

$$
\begin{equation*}
\tilde{\alpha}=\alpha D R \tag{137}
\end{equation*}
$$

### 2.4.2.1.1 Zenith angle computation

Figure 1 indicates the geometry involved in the computation of the zenith angle $\zeta$ at $P_{1}$.
Calculate:

$$
\begin{gather*}
\cos \tilde{\delta}=\sin \tilde{\varphi}_{1} \sin \tilde{\varphi}_{2}+\cos \tilde{\varphi}_{1} \cos \tilde{\varphi}_{2} \cos \left(\tilde{\lambda}_{2}-\tilde{\lambda}_{1}\right)  \tag{138}\\
\sin \tilde{\delta}=\sqrt{1-\cos ^{2} \tilde{\delta}}  \tag{139}\\
\tilde{\zeta}=\operatorname{atan} 2\left(\sin \tilde{\delta}, \cos \tilde{\delta}-\frac{r 1}{r 2}\right) \tag{140}
\end{gather*}
$$

being $\delta$ the Earth angle on the great circle connecting the receiver $\left(P_{1}\right)$ and the satellite $\left(P_{2}\right)$. The symbol $\operatorname{atan} 2(\mathrm{y}, \mathrm{x})$ indicates the function that computes the arctangent of $\mathrm{y} / \mathrm{x}$ with a range of $(-\pi, \pi]$.

FIGURE 1
Geometry of zenith angle computation


### 2.4.2.1.2 Ray-perigee computation

Figure 2 indicates the geometry involved in the computation of the coordinates of the ray-perigee $P_{P}$ : ray-perigee radius $r_{p}(\mathrm{~km})$, ray-perigee latitude $\varphi_{p}$ (degrees) and ray-perigee longitude $\lambda_{p}$ (degrees). Calculate $r_{p}$ :

$$
\begin{equation*}
r_{p}=r_{1} \sin \tilde{\zeta} \tag{141}
\end{equation*}
$$

Calculate $\varphi_{p}$ :
if $\left|\left|\varphi_{1}\right|-90^{0}\right|<10^{-10}$ use

$$
\varphi_{p}= \begin{cases}\zeta & \text { if } \varphi_{1}>0  \tag{142}\\ -\zeta & \text { if } \varphi_{1}<0\end{cases}
$$

otherwise use

$$
\begin{gather*}
\sin \tilde{\sigma}=\frac{\sin \left(\tilde{\lambda}_{2}-\tilde{\lambda}_{1}\right) \cos \tilde{\varphi}_{2}}{\sin \tilde{\delta}}  \tag{143}\\
\cos \tilde{\sigma}=\frac{\sin \tilde{\varphi}_{2}-\cos \tilde{\delta} \sin \tilde{\varphi}_{1}}{\sin \tilde{\delta} \cos \tilde{\varphi}_{1}}  \tag{144}\\
\tilde{\delta}_{p}=\frac{\pi}{2}-\tilde{\zeta}  \tag{145}\\
\sin \widetilde{\varphi}_{p}=\sin \widetilde{\varphi}_{1} \cos \widetilde{\delta}_{p}-\cos \widetilde{\varphi}_{1} \sin \widetilde{\delta}_{p} \cos \widetilde{\sigma}  \tag{146}\\
\cos \widetilde{\varphi}_{p}=\sqrt{1-\sin ^{2} \widetilde{\varphi}_{p}}  \tag{147}\\
\widetilde{\varphi}_{p}=\operatorname{atan} 2\left(\sin \widetilde{\varphi}_{p}, \cos \widetilde{\varphi}_{p}\right) \tag{148}
\end{gather*}
$$

Calculate $\lambda_{p}$ :
if $\left|\left|\varphi_{1}\right|-90^{\circ}\right|<10^{-10}$ use

$$
\varphi_{p}= \begin{cases}\zeta & \text { if } \varphi_{1}>0  \tag{149}\\ -\zeta & \text { if } \varphi_{1}<0\end{cases}
$$

otherwise use

$$
\begin{gather*}
\sin \left(\tilde{\lambda}_{1}-\tilde{\lambda}_{p}\right)=-\frac{\sin \widetilde{\sigma}_{\sin } \tilde{\delta}_{p}}{\cos \tilde{\varphi}_{p}}  \tag{150}\\
\cos \left(\tilde{\lambda}_{1}-\tilde{\lambda}_{p}\right)=\frac{\cos \widetilde{\delta}_{p}-\sin \tilde{\varphi}_{1} \sin \tilde{\varphi}_{p}}{\cos \tilde{\varphi}_{1} \cos \widetilde{\varphi}_{p}}  \tag{151}\\
\lambda_{p}=\left[\operatorname{atan} 2\left(\sin \left(\tilde{\lambda}_{1}-\tilde{\lambda}_{p}\right) \cos \left(\tilde{\lambda}_{1}-\tilde{\lambda}_{p}\right)\right)+\tilde{\lambda}_{1}\right] * R D \tag{152}
\end{gather*}
$$

being $\sigma$ the azimuth of $P_{2}$ seen from $P_{1}$ and $\delta_{p}$ the Earth angle between $P_{1}$ and the ray-perigee $P_{P}$.

## Geometry of ray-perigee computation



### 2.4.2.1.3 Great circle properties

Compute the great circle angle $\psi$ from ray-perigee to satellite:
if $\left|\left|\varphi_{p}\right|-90\right|<10^{-10}$ use

$$
\begin{equation*}
\Psi=\left|\varphi_{2}-\varphi_{p}\right| \tag{153}
\end{equation*}
$$

otherwise use

$$
\begin{gather*}
\cos \widetilde{\Psi}=\sin \widetilde{\varphi}_{p} \sin \widetilde{\varphi}_{2}+\cos \widetilde{\varphi}_{p} \cos \widetilde{\varphi}_{2} \cos \left(\tilde{\lambda}_{2}-\tilde{\lambda}_{p}\right)  \tag{154}\\
\sin \widetilde{\Psi}=\sqrt{1-\cos ^{2} \widetilde{\Psi}}  \tag{155}\\
\widetilde{\Psi}=\operatorname{atan} 2(\sin \widetilde{\Psi}, \cos \widetilde{\Psi}) \tag{156}
\end{gather*}
$$

Compute sine and cosine of azimuth $\sigma$ of satellite as seen from ray-perigee $\mathrm{P}_{\mathrm{p}}$ :
if $\left|\left|\varphi_{p}\right|-90\right|<10^{-10}$ use

$$
\begin{gather*}
\sin \tilde{\sigma}_{p}=0  \tag{157}\\
\cos \tilde{\sigma}_{p}\left\{\begin{array}{c}
-1 \text { if } \varphi_{p}>0 \\
1 \text { if } \varphi_{p}<0
\end{array}\right. \tag{158}
\end{gather*}
$$

otherwise use

$$
\begin{align*}
& \sin \widetilde{\sigma}_{p}=-\frac{\cos \widetilde{\varphi}_{2} \sin \left(\tilde{\lambda}_{2}-\tilde{\lambda}_{p}\right)}{\sin \widetilde{\Psi}}  \tag{159}\\
& \cos \widetilde{\sigma}_{p}=-\frac{\sin \widetilde{\varphi}_{2}-\sin \widetilde{\varphi}_{p} \cos \widetilde{\Psi}}{\cos \widetilde{\varphi}_{p} \sin \widetilde{\Psi}} \tag{160}
\end{align*}
$$

### 2.4.2.2 Integration endpoints

Indicating with $s_{1}$ and $s_{2}$ the distances of $P_{1}$ and $P_{2}$ respectively from the ray perigee compute:

$$
\begin{align*}
& s_{1}=\sqrt{r_{1}^{2}-r_{p}^{2}}  \tag{161}\\
& s_{2}=\sqrt{r_{2}^{2}-r_{p}^{2}} \tag{162}
\end{align*}
$$

### 2.4.2.3 Coordinates along the integration path: $c\left(h_{s} \varphi_{s}, \lambda_{s}\right)$

Being $s(\mathrm{~km})$ the distance of a point $P$ from the ray perigee $P_{P},\left(r_{p}, \varphi_{p}, \lambda_{p}\right)$ the ray perigee coordinates and $\sin \sigma_{p}, \cos \sigma_{p}$ the sine and cosine of the azimuth of the satellite as seen from the ray perigee, the coordinates of the point $P$ are calculated by the function $c$ as follows.
Inputs:
Distance $s(\mathrm{~km})$, ray-perigee coordinates $\left(r_{p}, \varphi_{p}, \lambda_{p}\right)$, sine and cosine of azimuth of satellite as seen from ray perigee $\sin \sigma_{p}, \cos \sigma_{p}$.

Outputs:
Coordinates of point $P: h_{s}(\mathrm{~km}), \varphi_{s}$ (degrees), $\lambda_{s}$ (degrees)
To compute the geocentric coordinates of any point $P$ (having distance $s$ from the ray perigee $P_{P}$ ) along the integration path, the following formulae have to be applied:

Calculate $h_{s}$ :

$$
\begin{equation*}
h_{s}=\sqrt{s^{2}-r_{p}^{2}}-R_{E} \tag{163}
\end{equation*}
$$

being $R_{E}$ the Earth mean radius.
Calculate great circle parameters:

$$
\begin{gather*}
\tan \tilde{\delta}_{s}=\frac{s}{r_{p}}  \tag{164}\\
\cos \widetilde{\delta}_{s}=\frac{1}{\sqrt{1+\tan ^{2} \widetilde{\delta}_{s}}}  \tag{165}\\
\sin \delta_{s}=\tan \widetilde{\delta}_{s} \cos \widetilde{\delta}_{s} \tag{166}
\end{gather*}
$$

Calculate $\varphi_{s}$ :

$$
\begin{gather*}
\sin \widetilde{\varphi}_{s}=\sin \widetilde{\varphi}_{p} \cos \widetilde{\delta}_{s}+\cos \widetilde{\varphi}_{p} \sin \widetilde{\delta}_{s} \cos \widetilde{\sigma}_{p}  \tag{167}\\
\cos \widetilde{\varphi}_{s}=\sqrt{1-\sin ^{2} \widetilde{\varphi}_{s}}  \tag{168}\\
\widetilde{\varphi}_{s}=\operatorname{atan} 2\left(\sin \widetilde{\varphi}_{s}, \cos \widetilde{\varphi}_{s}\right) * R D \tag{169}
\end{gather*}
$$

Calculate $\lambda_{s}$ :

$$
\begin{gather*}
\sin \left(\tilde{\lambda}_{s}-\tilde{\lambda}_{p}\right)=\sin \widetilde{\delta}_{s} \sin \widetilde{\sigma}_{p} \cos \widetilde{\varphi}_{p}  \tag{170}\\
\cos \left(\tilde{\lambda}_{s}-\tilde{\lambda}_{p}\right)=\cos \widetilde{\delta}_{s}-\sin \widetilde{\varphi}_{p} \sin \widetilde{\varphi}_{s}  \tag{171}\\
\lambda_{s}=\left[\operatorname{atan} 2\left(\sin \left(\tilde{\lambda}_{s}-\tilde{\lambda}_{p}\right), \cos \left(\tilde{\lambda}_{s}-\tilde{\lambda}_{p}\right)\right)+\tilde{\lambda}_{p}\right] * R D \tag{172}
\end{gather*}
$$

### 2.4.2.4 Slant TEC numerical integration

To compute slant TEC along a ray-path defined by its perigee coordinates, direction and end-point, a numerical integration algorithm is used. In NeQuick Gauss integration is used and is described in the following.

Inputs:

- $\quad h_{1}$, height of point $P_{1}(\mathrm{~km})$
- $\quad \varphi_{1}$, latitude of point $P_{1}$ (degrees)
- $\quad \lambda_{1}$, longitude of point $P_{1}$ (degrees)
- $\quad h_{2}$, height of point $P_{2}(\mathrm{~km})$
- $\quad \varphi_{2}$, latitude of point $P_{2}$ (degrees)
$-\quad \lambda_{2}$, longitude of point $P_{2}$ (degrees)
Output:
Slant TEC ( $10^{15} \mathrm{~m}^{-2}$ )
One possible numerical algorithm for slant TEC calculation is the following.
In the case of integration from ground to satellite ( $h_{1}<1000 \mathrm{~km}$ and $h_{2}>2000 \mathrm{~km}$ ) it is convenient to divide the integration path in three parts defining intermediate points $s_{a}, s_{b}$ :

$$
\begin{align*}
& s_{a}=\sqrt{\left(R_{E}+1000\right)^{2}-r_{p}^{2}}  \tag{173}\\
& s_{b}=\sqrt{\left(R_{E}+2000\right)^{2}-r_{p}^{2}} \tag{174}
\end{align*}
$$

We have that $\left(R_{E}+1000\right)^{2}=54334589.44$ and $\left(R_{E}+2000\right)^{2}=70076989.44$.
The slant TEC becomes therefore:

$$
\begin{equation*}
T E C=\int_{S_{1}}^{S_{a}} N(s) d s+\int_{S_{a}}^{S_{b}} N(s) d s+\int_{S_{b}}^{S_{2}} N(s) d s \tag{175}
\end{equation*}
$$

To compute each integral the algorithm described in $\S$ 2.4.2.5 can be used as:

$$
\begin{equation*}
\int_{g_{1}}^{g_{2}} N(s) d s=G N\left(g_{1}, g_{2}, \varepsilon, r_{p}, \sin \tilde{\varphi}_{p}, \cos \widetilde{\varphi}_{p}, \sin \tilde{\sigma}_{p}, \cos \tilde{\sigma}_{p}, \lambda_{p}\right) \tag{176}
\end{equation*}
$$

where the parameter $\varepsilon$ indicates the integration accuracy. Here it is assumed that:

- $\quad \varepsilon=0.001$ for the integration between $s_{1}$ and $s_{a}$
- $\quad \varepsilon=0.01$ for the integrations between $s_{a}$ and $s_{b}$ and between $s_{b}$ and $s_{2}$


### 2.4.2.5 Gauss algorithm

Inputs:

- $\quad$ distances from the ray perigee of the first integration endpoint: $g_{1}(\mathrm{~km})$
- distances from the ray perigee of the second integration endpoint: $g_{2}(\mathrm{~km})$
- Integration accuracy $\varepsilon$
- Ray-perigee parameters: $r_{p}(\mathrm{~km}), \sin \widetilde{\varphi}_{p}, \cos \widetilde{\varphi}_{p}, \sin \widetilde{\sigma}_{p}, \cos \widetilde{\sigma}_{p}, \lambda_{\mathrm{p}}($ degrees $)$

Output:
TEC [TECu]

To be able to compute NeQuick electron density, in all the following computations it is necessary to calculate the coordinates of the point $P$ along the ray-path using the algorithm illustrated in § 2.4.2.3.

$$
\begin{equation*}
(h(s), \varphi(s), \lambda(s))=c\left(s, \sin \widetilde{\varphi}_{p}, \cos \widetilde{\varphi}_{p}, \sin \widetilde{\sigma}_{p}, \cos \widetilde{\sigma}_{p}, \tilde{\lambda}_{p}\right) \tag{177}
\end{equation*}
$$

Start the calculation using 8 points:

$$
\begin{equation*}
n=8 \tag{178}
\end{equation*}
$$

Repeat the following computations until the desired precision is obtained.
Calculate the integration intervals:

$$
\begin{gather*}
\Delta n=\frac{g_{2}-g_{1}}{n}  \tag{179}\\
g=0.5773502691896 \cdot \Delta n  \tag{180}\\
y=g_{1}+\frac{\Delta n-g}{2}  \tag{181}\\
G N_{2}=\frac{\Delta n}{2} \sum_{i=0}^{n-1}\left[f\left(y+i \Delta_{n}\right)+f\left(y+i \Delta_{n}+g\right)\right] \tag{182}
\end{gather*}
$$

Double the number of points:

$$
\begin{equation*}
n=2 n \tag{183}
\end{equation*}
$$

and define

$$
\begin{equation*}
G N_{1}=G N_{2} \tag{184}
\end{equation*}
$$

repeating steps from (178) to (181) it is now possible to compare the two values obtained to see if the required accuracy is achieved: if

$$
\begin{equation*}
\left|G N_{1}-G N_{2}\right|>\varepsilon\left|G N_{1}\right| \tag{185}
\end{equation*}
$$

then continue increasing the number of points (183), redefine $G N_{1}$ (184) and repeat again steps from (178) to (182).

When the test (185) fails, the required accuracy has been reached, and the value of the integral is obtained by:

$$
\begin{equation*}
T E C=\left(G N_{2}+\frac{G N_{2}-G N_{1}}{15}\right) \times 10^{-13} \tag{186}
\end{equation*}
$$

### 2.5 Changes introduced in NeQuick2

The changes introduced in NeQuick 2.1 with respect to the first version of the model are summarized. The most important modifications introduced are related to the bottomside electron density profile representation and to the shape parameters of the topside ionosphere.

### 2.5.1 Bottomside modification

The F1 layer formulation has been reformulated [Leitinger et al., 2005, RD2] to avoid the occurrence of the anomalous behaviour of the model reported at heights in the E and F1 layer of the ionosphere.

As listed in § 2.2.4, here the new parameters are provided as:

$$
\begin{gather*}
h m F 1=\frac{h m E+h m F 2}{2}  \tag{187}\\
f o F 1=\left\{\begin{array}{cc}
0 & \text { if } \quad f o E<2 \\
1.4 f o E & \text { if } \quad f o E \geq 2 \text { and } 1.4 f o E \leq 0.85 f o F 2 \\
0.85 \cdot 1.4 f o E & \text { if } 1.4 f o E>0.85 f o F 2
\end{array}\right.  \tag{188}\\
B 1_{\text {top }}=0.3(h m F 2-h m F 1)  \tag{189}\\
B 1_{b o t}=0.5(h m F 1-h m E)  \tag{190}\\
\mathrm{BE}_{\text {top }}=\max (0.5(h m F 1-h m E), 7) \tag{191}
\end{gather*}
$$

$f o E, f o F 1$ and $f o F 2$ are given in $(\mathrm{MHz}) ; h m E, h m F 1$ and $h m F 2$ are given in (km).

### 2.5.2 Topside modification

The topside of the $F 2$ layer of the ionosphere is implemented using a semi-Epstein layer representation. The electron density as function of height $N(h)$ in the topside is represented using a height dependent thickness parameter $H_{0}$ [RD3]:

$$
\begin{gather*}
N(h)=\frac{4 N m F 2}{(1+\exp (z))^{2}} \exp (z)  \tag{192}\\
z=\frac{h-h m F 2}{H_{0}\left[1+\frac{r g(h-h m F 2)}{r H_{0}+g(h-h m F 2)}\right]} \tag{193}
\end{gather*}
$$

in which the constant parameters $r=100$ and $g=0.125$ are used. NeQuick 2.1 includes a simplification of the topside thickness parameter $H_{0}$, using a single empirical relation to represent its correlation to the $F 2$ bottomside thickness parameter $B 2_{\text {bot }}$ [Nava et al., 2008 RD1]:

$$
\begin{gather*}
H_{0}=k B 2_{\text {bot }}  \tag{194}\\
k=3.22-0.0538 f o F 2-0.00664 h m F 2+0.113 \frac{h m F 2}{B 2_{\text {bot }}}+0.00257 R_{12} \tag{195}
\end{gather*}
$$

with a value of 1 imposed on $k$ as a lower limit.

### 2.5.3 Modip map

To get the magnetic coordinates needed by the model for computing the F2 layer peak and F1 layer height the NeQuick ITU-R (version 1) used a map of dip-latitude (diplats.asc). It was necessary to compute both Modip and Dip at the location of interest. The change in the formulation of the F1 layer in the NeQuick 2.1 now requires only the computation of the Modip.
NeQuick 2.1 model uses a grid of Modip values contained in the file modip.asc. To compute the Modip value in a point at latitude $\varphi$ and longitude $\lambda$, a third order interpolation is applied as it is explained in $\S 2.3 .3 .1$. The Modip map included in the package of version 2.1 has been generated with the magfit subroutine. The magfit subroutine is patterned after the Fortran subroutine in the ITS propagation package, 2006, called iongrid, provided by ESA and available at ITU-R in support of Recommendation ITU-R P.1239. It calculates the magnetic dip angle and the gyrofrequency for the
given reflection height and location. This calculation is described in § 2 of Recommendation ITU-R P.1239-3 (02-2012).

### 2.5.4 Other package and source code modifications

The set of CCIR coefficients that have been used in previous versions of the model have been replaced by the set of ITU-R coefficients provided in the iongrid program (software provided by ESA and available at ITU-R in support of Recommendation ITU-R P.1239).
From the point of view of FORTRAN programming, version 2 of NeQuick includes various improvements. In particular to avoid the use of the ENTRY statement, new subroutines have been included: prepNeQ and NeNeQ.

Version 2.1 corrects some formal bugs of earlier NeQuick versions in the management of modip maps. Some computational improvements have also been introduced.

### 2.6 References

[RD1] Nava, B., Coisson, P., Radicella, S.M., "A new version of the NeQuick ionosphere electron density model", J.of Atmospheric and Solar-Terrestrial Physics, Iss. 70, 1856-1862, 2008.
[RD2] Leitinger, R., Zhang, ML, Radicella, S/M., "An improved bottomside for the ionospheric electron density model NeQuick", Annals of Geophysics, 48(3), p. 525-534, 2005.
[RD3] Coisson, P., Radicella, S.M, Leitinger, R., Nava, B., "Topside electron density in IRI and NeQuick: features and limitations", Adv. Space Research., 37 (5), p. 937-942, 2006.
[RD4] Rawer K., "Propagation of decameter waves (HF-band)", in B. Landmark (Ed), Meteorological and astronomical influences on radio wave propagation, New York Academic Press, p. 221-250, 1963.

## 3 Vertical total electron content maps

A set of grid point maps of Vertical Total Electron Content (VTEC) representative of various levels of solar activity are considered of interest for model verification or preliminary assessment of various trans-ionospheric propagation effects such as Faraday rotation, propagation delay, refraction, angle of arrival variation among others. They may be of interest for evaluation of HF propagation. Those maps are available from the Radiocommunication Bureau.

A de facto standard widely used by various scientific communities is proposed for the format of the provided VTEC maps. Such format is the IONEX (IONosphere map EXchange) format, see details in: http://igscb.jpl.nasa.gov/igscb/data/format/ionex1.pdf.
The International GNSS Service (IGS, http://igscb.jpl.nasa.gov/components/prods.html) routinely generates Global Ionospheric Maps (GIMs) in IONEX format on a grid resolution of $2^{\mathrm{h}} \times 5^{\circ} \times 2.5^{\circ}$ in UT, longitude and latitude for one day GIMs are routinely generated by four processing centres using dual-frequency GPS observations on a large network of permanent reference stations around the globe, and a final product is obtained by combining the maps from all four processing centres.
Even if monthly median or mean maps together with its inter-quartile ranger or its standard deviation are provided, the next question is whether they are sufficiently representative of the climatology of the ionosphere. An example of the day-to-day variability and median for 30 days in April 2000 is presented in Fig. 3. In such an example, it is observed that the median curve does not clearly show the very variable equatorial anomaly crests seen in a number of individual days. Another question is related to the symmetry and skewness of the daily values. Examples of the problems of mean/median values and variability are further discussed in the attached presentation.

FIGURE 3
Meridional cross-section of Vertical TEC from IGS combined maps, April 2000 at 13:00 UT, Day 1 to 30 and median (thick line)


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While the problem of the selection of what can be considered representative is not solved, it is still considered of practical use to incorporate in ITU-R databanks a set of monthly mean maps derived from observations in an accepted grid format, as examples, always remarking the limitations of the averaging and the variability of the ionosphere.
The following test cases, have been derived:

- $\quad$ Three year related to three types solar activity:
- Solar maximum (2002)
- Solar minimum (2009)
- Solar intermediate activity (2005)
- Four sets of 31 days corresponding to different annual seasons, they are derived as 15 days plus and minus the following days:
- Day of year 79 (Northern hemisphere spring equinox)
- Day of year 171 (Northern hemisphere summer solstice)
- Day of year 265 (Northern hemisphere autumn equinox)
- Day of year 350 (Northern hemisphere winter solstice)

IONEX format supports the inclusion of the VTEC and its standard deviation error. In the VTEC values, the mean value for each grid point over the 31 days has been extracted, and in the standard deviation error, the standard deviation for the 31 days is given.

A set of figures showing the VTEC map and its standard deviation for the three solar activity cases and the month around the Northern hemisphere spring equinox are presented in Fig. 4 as examples of absolute values and variability.

In the future, in order to incorporate information about the range of variability observed around mean values, in addition to the mean maps and its standard deviation, it is suggested to incorporate three daily maps based on certain criteria such as:

- The daily map which minimizes in all points and hours the root mean square error to the mean value.
- The daily map with maximum positive bias (difference between daily map and mean map for all points and hours).
- The daily map with maximum negative bias (difference between daily map and mean map for all points and hours).


## FIGURE 4

Mean (left) and standard deviation (right) of VTEC daily maps for a month around the Northern hemisphere spring equinox for low (top), intermediate (middle) and high (bottom) solar activity at 14:00 UT (15:00 UT for high solar activity)


