#### Documentation of the GKW to GKDB data conversion

Y. Camenen on behalf of the GKDB working group

Last update: February 9, 2018

### Preface

This document describes how to transform inputs and outputs from a GKW flux-tube simulation to match the format used in the GyroKinetic DataBase (GKDB). The reader is assumed to have some knowledge of GKW and to have read the documentation of the GKDB.

### Contents

1	Preface	1
2	Conventions and normalisations 2.1 Coordinate systems	<b>3</b>
	2.2 Reference quantities	4
3	Inputs	5
	3.1 Magnetic equilibrium	5
	3.1.1 Plasma shape	5
	3.1.2 Poloidal angle	5
4	Outputs	6

#### Conventions and normalisations

#### 2.1 Coordinate systems

In GKW, the toroidal direction is defined to have the cylindrical coordinate system  $(R, Z, \varphi)$  right-handed whereas in the GKDB it is defined to have  $(R, \varphi, Z)$  right-handed, see Fig.2.1. In practice, it means that:

$$\varphi^{\text{GKW}} = -\varphi^{\text{GKDB}} \tag{2.1}$$

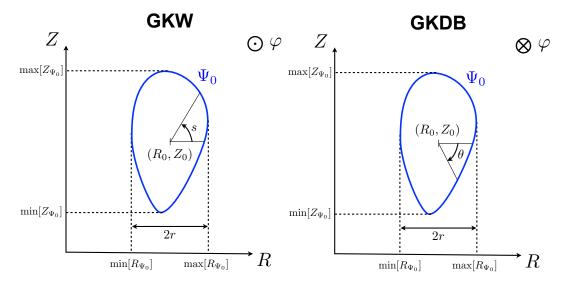


Figure 2.1: Cylindrical coordinate system used in GKW (left) and the GKDB (right).

The flux surface centre definition depends on how the magnetic equilibrium is specified. For miller geometry, the definition of  $R_0$  is identical to that used in the GKDB and  $Z_0$  is given as an input in the geometry namelist:

$$R_0^{\text{GKW-miller}} = R_0^{\text{GKDB}}$$
  $Z_0^{\text{GKW-miller}} = \text{zmil}R_{\text{ref}}^{\text{GKW}}$  (2.2)

For chease geometry,  $R_0$  is taken to be the value of ROEXP specified in the hamada.dat file and  $Z_0$  is the elevation of the magnetic axis.

$$R_0^{\rm GKW-chease} = {\rm ROEXP} \hspace{1cm} Z_0^{\rm GKW-chease} = Z_{\rm axis} \hspace{1cm} (2.3)$$

The definition of the (dimensional) radial coordinate r is identical in GKW and the GKDB:

$$r^{\text{GKW}} = r^{\text{GKDB}} \tag{2.4}$$

The calculation of the poloidal angle  $\theta$  used in the GKDB from GKW inputs is documented in section 3.1. At this stage, just notice that most of the time  $Z_0^{\rm GKW} \neq Z_0^{\rm GKDB}$ , therefore, the points s=0 and  $\theta=0$ do not necessarily coincide.

#### 2.2 Reference quantities

The conversion from GKW to GKDB requires to know the ratio of the reference quantities used in both formats. These ratio are:

$$\begin{split} q_{\mathrm{rat}} &= \frac{q_{\mathrm{ref}}^{\mathrm{GKW}}}{q_{\mathrm{ref}}^{\mathrm{GKDB}}} = -\frac{1}{\mathbf{z}_{e^{-}}^{\mathrm{GKW}}} & L_{\mathrm{rat}}^{\mathrm{miller}} &= \frac{R_{\mathrm{ref}}^{\mathrm{GKW-miller}}}{R_{\mathrm{ref}}^{\mathrm{GKDB}}} = 1 \\ m_{\mathrm{rat}} &= \frac{m_{\mathrm{ref}}^{\mathrm{GKW}}}{m_{\mathrm{ref}}^{\mathrm{GKDB}}} = \frac{m_e}{m_D} \frac{1}{\mathrm{mass}_{e^{-}}^{\mathrm{GKW}}} & B_{\mathrm{rat}}^{\mathrm{miller}} &= \frac{B_{\mathrm{ref}}^{\mathrm{GKW-miller}}}{B_{\mathrm{ref}}^{\mathrm{GKDB}}} = 1 \\ T_{\mathrm{rat}} &= \frac{T_{\mathrm{ref}}^{\mathrm{GKW}}}{T_{\mathrm{ref}}^{\mathrm{GKDB}}} = \frac{1}{\mathrm{temp}_{e^{-}}^{\mathrm{GKW}}} & L_{\mathrm{rat}}^{\mathrm{chease}} &= \frac{R_{\mathrm{ref}}^{\mathrm{GKW-chease}}}{R_{\mathrm{ref}}^{\mathrm{GKDB}}} = \frac{R_{\mathrm{OEXP}}^{\mathrm{EXP}}}{R_{\mathrm{OKDB}}^{\mathrm{GKDB}}} \\ n_{\mathrm{rat}} &= \frac{n_{\mathrm{ref}}^{\mathrm{GKW}}}{n_{\mathrm{ref}}^{\mathrm{GKDB}}} = \frac{1}{\mathrm{dens}_{e^{-}}^{\mathrm{GKW}}} & B_{\mathrm{rat}}^{\mathrm{chease}} &= \frac{B_{\mathrm{ref}}^{\mathrm{GKD-chease}}}{B_{\mathrm{ref}}^{\mathrm{GKDB}}} = \frac{B_{\mathrm{OEXP}}}{B_{\mathrm{OKDB}}^{\mathrm{GKDB}}} \end{split}$$

The  $e^-$  subscript denotes the electron species and the electron to deuterium mass ratio is taken to be  $\frac{m_e}{m_D} = 2.7237 \times 10^{-4}$  in the GKDB.

The poloidal asymmetry factor for the density can be computed from data in the cfdens file.

With chease geometry, the ratios  $L_{\text{rat}}$  and  $B_{\text{rat}}$  can be computed from data in the hamada.dat file.

### Inputs

#### 3.1 Magnetic equilibrium

#### 3.1.1 Plasma shape

Let's call  $\{R_{\Psi_0}, Z_{\Psi_0}\}$  a set of points describing the flux surface of interest. Only miller and chease magnetic equilibrium specifications are compatible with the GKDB format.

#### 3.1.2 Poloidal angle

# Outputs

## Bibliography