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# Multiscale modeling of ice deformation behavior

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# Abstract

Understanding the flow of ice in glaciers and polar ice sheets is of increasing relevance in a time of potentially significant climate change. The flow of ice has hitherto received relatively little attention from the structural geological community. This paper aims to provide an overview of methods and results of ice deformation modeling from the single crystal to the polycrystal scale, and beyond to the scale of polar ice sheets. All through these scales, various models have been developed to understand, describe and predict the processes that operate during deformation of ice, with the aim to correctly represent ice rheology and self-induced anisotropy. Most of the modeling

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tools presented in this paper originate from the material science community, and are currently used and further developed for other materials and environments. We will show that this community has deeply integrated ice as a very useful "model" material to develop and validate approaches in conditions of a highly anisotropic behavior. This review, by no means exhaustive, aims at providing an overview of methods at different scales and levels of complexity.

Keywords: ice mechanical behavior, multiscale modeling, viscoplastic anisotropy, fabric development

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#### 1. Introduction

Ice is a common mineral on the Earth's surface, where it occurs as ice
Ih. As ice is relatively close to its melting temperature, glaciers and polar
ice sheets deform by ductile dislocation creep at strain rates in the order
of 10<sup>-12</sup> to 10<sup>-6</sup> s<sup>-1</sup>. Research on the flow of ice is of direct importance
to society as it is needed to understand and predict the effects that global
warming could have on sea level rise, glacier retreat, etc. There is also an
increasing awareness that ice is a valuable analogue for other minerals and
crystalline materials, as it is the only common mineral where this creep can
be readily observed in nature and in the laboratory. Numerical modeling
has become a key method to link the mechanics of ice from the dislocation
scale to that of flowing ice masses.

Most of the efforts made to simulate the ductile mechanical behavior of polycrystalline ice are related to the modeling of ice flow and fabric evolution in the conditions of polar ice sheets or glaciers. Ice is increasingly considered a model material to validate micro-macro mechanical approaches for materials with a high viscoplastic anisotropy. Most of the modeling techniques presented in this paper are currently used or further developed for other materials. For geological applications, one main limitation could be related to the "one phase" approach for most of these techniques, well adapted to ice. The reader will find, at the end of the paper, a table summarizing the main aspects of each techniques, with application ranges and limitations.

#### 1.1. Mechanical properties of ductile ice

Ice Ih has an hexagonal crystal structure with a c/a ratio of 1.628. This 64 c/a ratio is very close to the 1.633 value for a closely packed structure, but ice is not closely packed (see Schulson and Duval (2009) for a recent review). The elastic anisotropy of ice single crystals is small. The Young modulus E67 only varies by about 30%, depending on the direction of the loading axis with respect to the c-axis. The highest value is along the c-axis with E = 11.8GPa at -16°C (Gammon et al., 1983). Single crystals deform plastically essentially by glide of dislocations on 71 the basal plane. There are three equivalent  $< 1\overline{2}10 >$  directions for the Burgers vector, but slip on the basal plane is almost isotropic. In conditions where basal slip is favored, the stress-strain rate relationship after a strain of about 5\% can be expressed by a power law with a stress exponent  $n=2\pm0.3$ (Higashi et al., 1965; Jones and Glen, 1969; Mellor and Testa, 1969). At similar strain rates, the equivalent stress requested for non-basal slip is about 77 60 times larger than for basal slip (Duval et al., 1983). For ice polycrystals deformed under the laboratory conditions (strain rate between about  $10^{-8}$  s<sup>-1</sup> and  $10^{-6}$  s<sup>-1</sup> and temperature generally higher than -30°C), strain is essentially due to intracrystalline dislocation glide. The transient creep regime is characterized by a strong directional hardening until the strain-rate minimum is reached for an overall strain of 1% (Duval et al., 1983). This strain-rate decrease can reach three orders of magnitude. 84 It is associated to the development of a strong internal stress field due to plastic incompatibility between grains (Ashby and Duval, 1985; Duval et al., 1983; Castelnau et al., 2008b). A significant part of the transient creep is recoverable, i.e., on unloading a creep specimen, a reverse creep is observed,

with reverse strain which can be more than ten times the initial elastic strain

Ouval, 1976; Duval et al., 1983). In the secondary creep regime, isotropic polycrystals deform (at similar stress levels) a 100 times slower than a single crystal optimally oriented for basal slip. In this regime, the minimum strain rate and the stress are linked by a power law, referred to as Glen's law in glaciology (Glen, 1955), expressed through a relationship of the form (1) for temperatures lower that -10°C.

with  $\bar{\sigma}$  the applied stress,  $E_p=0.72$  eV and the stress exponent n=3

$$\dot{\bar{\varepsilon}}_{min} = A\bar{\sigma}^n exp(-E_p/k_B T) \tag{1}$$

(Barnes et al., 1971; Budd and Jacka, 1989). A is a constant,  $k_B$  the Boltz-97 mann constant and T the temperature. Above -10°C,  $\dot{\bar{\varepsilon}}_{min}$  rises more rapidly 98 with increasing temperature and cannot be described by this equation (Mor-99 gan, 1991). No grain-size effect is expected for power-law secondary creep 100 at laboratory conditions (see Duval and Le Gac (1980); Jacka (1994) for 101 instance). But a grain size effect was, however, measured during transient 102 creep (Duval and Le Gac, 1980). 103 At strains larger than 1 to 2\% (tertiary creep regime), dynamic recrystalliza-104 tion is predominant, and new grain microstructures and crystal orientations 105 are generated (Jacka and Maccagnan, 1984; Duval et al., 2000). 106 At stresses lower than 0.1 MPa, relevant to deformation conditions in glaciers, 107 ice sheets or planetary bodies, there is a clear indication of a creep regime 108 with a stress exponent lower than two. This indication results from both 109 the analysis of field data and laboratory tests, although the difficulty of 110 obtaining reliable data at strain rates lower than  $10^{-10}$ s<sup>-1</sup> is at the ori-111 gin of contradictory results (Mellor and Testa, 1969; Barnes et al., 1971; 112 Dahl-Jensen and Gundestrup, 1987; Pimienta et al., 1987; Lipenkov et al., 1997; Goldsby and Kohlstedt, 1997). In particular, Goldsby and Kohlstedt

(1997) suggest a grain-size dependence of the ice viscosity associated with 115 this low stress regime, based on laboratory experiments performed on very 116 small grain-size samples. This grain-size effect would be associated with a grain boundary-sliding dominated creep. Its extrapolation to polar ice-118 core deformation conditions remains controversial (Duval and Montagnat, 119 2002). Diffusional creep, commonly associated with such conditions in many 120 materials yields a viscosity much higher than that deduced from field data 121 (Lliboutry and Duval, 1985). For a review on ice behavior, see (Duval et al., 122 2010). 123 Ice as a model material exhibits a challenging viscoplastic anisotropy ow-124 ing to the presence of only two independent easy slip systems for the dis-125 locations (basal plane). While five independent systems are required to 126 accommodate an arbitrary deformation in a single crystal (Taylor, 1938), 127 Hutchinson (1977) showed that four systems are required for allowing an 128 hexagonal polycrystal such as ice to deform. Being able to represent and to 129 take into account this anisotropy in micro-macro models which aim at link-130 ing the single crystal scale to the polycrystal scale, is of primary interest to 131 the material science community. This anisotropy needs to be accounted for 132 at the dislocation scale in order to build physically-based model for the acti-133 vation of (poorly known) secondary slip systems. The impact of dislocation 134 induced internal stress fields, but also the characterization and development 135 of highly heterogeneous strain and stress fields within polycrystals, and their 136 impact on fabric development turn out to be of strong importance (Castel-137 nau et al., 1996a; de la Chapelle et al., 1998). 138 During gravity-driven flow of glaciers and ice sheets, the macroscopic 139 behavior of ice becomes progressively anisotropic with the development of fabrics (or textures, c-axis preferred orientations). This anisotropy and its

development depends on the flow conditions, but strongly influences the response of ice layers to imposed stress (see Gundestrup and Hansen (1984); 143 Van der Veen and Whillans (1990); Mangeney et al. (1997) for pioneer field work and modeling on the subject). Indeed, a polycrystal of ice with most 145 of its c-axes oriented in the same direction deforms at least ten times faster 146 than an isotropic polycrystal, when sheared parallel to the basal planes. Fabrics basically develop as the result of lattice rotation by intracrystalline 148 slip (Azuma and Higashi, 1985; Alley, 1988, 1992). Dynamic recrystalliza-149 tion can have a major impact on fabric development, especially at temper-150 atures above -10°C close to bedrocks or within temperate glaciers (Alley, 151 1992; Duval and Castelnau, 1995; de la Chapelle et al., 1998; Montagnat 152 et al., 2009), see Section 5. Questions, however, remain to what extent dif-153 ferent recrystallization processes operate as a function of depth in polar ice sheets (Kipfstuhl et al., 2006, 2009; Weikusat et al., 2009). 155

#### 156 1.2. Main objectives

Accurate modeling of ice flow under natural conditions is relevant for 157 many scientific objectives, such as the response of ice sheet to climate 158 changes (Seddik et al., 2012), the interpretation of climate signals extracted 159 from ice cores (Faria et al., 2010), the energy balance in extraterrestrial 160 satellites (Sotin et al., 2009), and since a few years, the accurate prediction 161 of sea-level rise that is linked to the behavior of fast-moving coastal glaciers 162 (Gillet and Durand, 2010). In this context, challenges are mainly (i) to 163 establish an ice flow law adapted to low stress conditions, changes in tem-164 peratures and impurity content, (ii) to consider the macroscopic anisotropy 165 due to fabric development at the given conditions, (iii) to be able to inte-166 grate processes such as dynamic recrystallization that can strongly influence 167

fabric development and the flow law.

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The aim of this paper is to present a general overview of the main modeling techniques adapted to ice, and the main modeling results obtained from the single crystal scale to the large scale that is relevant to ice sheet flow modeling. Techniques are highly diverse, from dislocation dynamics (micron scale) to Finite Element methods that are adapted to the whole ice sheet (km scale), via mean-field and full-field micro-macro approaches and coupling with a microstructure evolution models (cm to m scale, limited to a 2D configuration, see 5.2). We will mostly focus on recent advances and topics that are still under development.

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#### 2. Modeling ice single crystal behavior

Owing to its high viscoplastic anisotropy, with dislocations gliding mostly 181 on the basal plane, studying and modeling ice single crystal behavior is a 182 challenge for regular approaches. 183 Recent efforts focused on three main objectives; (i) understanding, repre-184 senting and taking into account the dislocation dynamics, (ii) improving our 185 knowledge about secondary slip systems in ice, (iii) providing an accurate 186 crystal plasticity constitutive law that can be implemented in mean-field and full-field approaches for micro-macro polycrystal models. For the two first 188 objectives, Dislocation Dynamic models (DD) were used at the scale of the 189 interaction between dislocation populations (Section 2.1). At a larger scale, the Field Dislocation Mechanic modeling approach (FDM) was applied to 191 ice to evaluate the role of internal stresses associated with dislocation fields

and arrangements (Section 2.2). Section 2.3 presents a crystal-plasticity model adapted to the transient creep behavior of ice single crystals.

# 2.1. Dislocation Dynamics modeling

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Dislocation dynamics in ice was shown to be scale free and intermit-196 tent, thanks to dislocation avalanche measurements via acoustic emissions 197 (Weiss and Grasso, 1997; Weiss et al., 2001; Weiss and Marsan, 2003; Weiss 198 and Montagnat, 2007). Ice was used as a model material for the following 199 reasons: (i) transparency allows direct verification that acoustic emission 200 activity is not related to microcraking, (ii) with the range of stress and 201 temperature considered, diffusion creep is not a significant mechanism, and 202 deformation occurs by dislocation glide only. DD modeling tools were used 203 to better understand and characterize this scale free and intermittent be-204 havior (for example Miguel et al. (2001); Weiss and Miguel (2004)). 205 Miguel et al. (2001) made use of a discrete dislocation dynamics model with 206 a two-dimensional cross-section of the crystal. This 2D space is randomly 207 filled with edge dislocations gliding along a single slip direction parallel to 208 their respective Burgers vector. This simplification is an effective way to 209 describe materials like ice crystals owing to their strong plastic anisotropy 210 with a single slip system dominating. A basic feature common to most DD 211 models is that dislocations interact with each other through the long-range 212 elastic stress field they produce in the host material. In (Miguel et al., 213 2001), dislocation velocity depends linearly on this effective stress, and the 214 Peierls stress is set to zero. Mechanisms for dislocation annihilation and 215 multiplication are classically taken into account. 216

Within this simplified scheme the authors found that dislocations generate a slowly evolving configuration landscape which coexists with rapid col-

lective rearrangements. These arrangements involve a comparatively small fraction of dislocations and lead to an intermittent behavior of the net plastic 220 response. The model was therefore able to reproduce the fact that disloca-221 tions themselves, through the various structures such as dipoles and walls, 222 generate a pinning force landscape that is virtually frozen into a slow state. 223 Creation and annihilation mechanisms allow the system to jump between 224 slow dynamics states through bursts of activity. 225 More recently, Chevy et al. (2007, 2012) used DD simulations to analyze 226 torsion tests performed on ice single crystals. The tests were performed 227 with the ice-crystal c-axis oriented parallel to the torsion axis so that basal 228 screw dislocations were mainly activated. With synchrotron topography 229 analyses of the deformed samples, it was possible to show that dislocation 230 arrangements were highly heterogeneous, with a scale-invariant character 231 and long-range correlations (Montagnat et al., 2006; Weiss and Montagnat, 232 2007; Chevy et al., 2010). Although these tests were performed in a way that 233 highly favored basal glide, the double-cross slip mechanisms was invoked to 234 explain this scale invariant dislocation arrangement. 235 Three-dimensional DD simulations, based on the TRIDIS code (Verdier 236 et al., 1998), were adapted to these torsion tests on ice and the hexago-237 nal structure. Screw dislocation sources were positioned within one slip 238 plane at the periphery of a cylinder submitted to a constant torque. Cross-239 slip on prismatic planes was made possible thanks to the internal stress 240 induced by the pile-up of basal dislocations in the center of the cylinder 241 (where  $\sigma_{app} = 0$ ), which produces the out-of-plane component needed (see 242 Fig. 1). Simulation results allowed to test this hypothesis, and explain the power law relationship between stress and strain rate (Chevy et al., 2012).

## 2.2. Field Dislocation Mechanics (FDM)

Field dislocation theory is a mesoscale approach, which aims at tak-246 ing into account the inhomogeneous distribution of dislocations in plasticity modeling. Therefore, FDM modeling makes it possible to represent and con-248 sider the internal stress field created by the dislocation arrangements within 249 the crystal. FDM is a continuous approach able to deal simultaneously with 250 long-range correlations associated with distortion fields, internal stresses due 251 to dislocation arrangements, and short-range correlations (Acharya, 2001). 252 The reader is referred to (Acharya and Roy, 2006; Varadhan et al., 2006; 253 Fressengeas, 2010) for details. 254 The first application to ice samples was performed in the configuration of 255 the torsion test presented in part 2.1 (Taupin et al., 2007). This test is 256 by itself highly heterogeneous, and this heterogeneity was shown to induce 257 unexpected non-basal slip. Taking into account the coupled dynamics of 258 geometrically necessary screw dislocations gliding in the basal plane (also 259 called "excess" dislocations) and statistical dislocations developed through 260 cross slip in prismatic planes, the model was able to reproduce the creep 261 curves during torsion, and the size effect measured experimentally (see Fig. 262 2). More recently, the model was used to reproduce the complex scale-263 invariant character of dislocation arrangements forming during torsion tests 264 on ice single crystals (Chevy et al., 2010). In particular, the fact that the 265 model takes into account both the long-range elastic interactions due to 266 the presence of dislocations and the short-range interactions inherent to the 267 transport of dislocations (obstacles, cross-slip, etc.) allowed to reproduce 268 the shift in control of the dislocation distribution by long-range correlations 269 at low strain to a control by short-range correlations at strain as high as 270 50%. It was shown that non-basal dislocations activated by the internal stress fields induce a screening potential at large strain, through obstacles such as twist sub-boundaries. However, this screening was shown to be too small to hinder creep acceleration prevailing during torsion creep test on ice single crystals (Chevy et al., 2010).

# 2.3. Crystal plasticity modeling

Constitutive relations to describe the transient creep of ice single crystals
have been proposed by Castelnau et al. (2008b) and then used in a modified
version in Suquet et al. (2011). One of the difficulty here is the description
of the softening of basal slip in the transient regime, as discussed above.
As is usual in crystal plasticity at infinitesimal strains, the strain tensor is
decomposed into the sum of an elastic  $\varepsilon^{e}$  and a viscoplastic  $\varepsilon^{vp}$  part

$$\varepsilon = \varepsilon^{\mathsf{e}} + \varepsilon^{\mathsf{vp}}$$
 . (2)

The elastic strain is related to the local stress tensor  $\sigma$  with the local compliance tensor S, and the viscoplastic strain results from slips on a total of M different slip systems:

$$\varepsilon^{\mathsf{e}} = \mathbf{S} : \boldsymbol{\sigma}, \quad \varepsilon^{\mathsf{vp}} = \sum_{k=1}^{M} \gamma^{(k)} \boldsymbol{\mu}^{(k)}.$$
(3)

Here,  $\boldsymbol{\mu}^{(k)} = \frac{1}{2}(\mathbf{n}^{(k)} \otimes \mathbf{b}^{(k)} + \mathbf{b}^{(k)} \otimes \mathbf{n}^{(k)})$  is the (purely geometric) Schmid tensor depending on the orientation of the slip system (k),  $\mathbf{n}$  being the slip plane normal and  $\mathbf{b}$  the slip direction (parallel to the Burgers vector and orthogonal to  $\mathbf{n}$ ) in that plane, with  $\otimes$  the dyadic product. Ice crystals, which have an hexagonal symmetry, deform easily by shear on the basal plane, on the three systems  $\{0001\} < 11\overline{2}0 >$  which provide only two independent systems. The three prismatic systems  $\{1\overline{1}00\} < 11\overline{2}0 >$ provide two more independent systems. An additional independent slip

system is thus required to attain any isochoric deformation at the single 294 crystal level and this is achieved by adding the six  $\langle c + a \rangle$  pyramidal 295 systems  $\{11\overline{2}2\} < 11\overline{2}3 >$ . In total, M = 12 slip systems are taken into 296 account in the present analysis. 297 In the constitutive relations originally proposed by Castelnau et al. 298 (2008b), the slip rate on the k-th system is related to the resolved shear stress  $\tau^{(k)}$  on that system through:

$$\dot{\gamma}^{(k)} = \dot{\gamma}_0^{(k)} \left( \frac{\left| \tau^{(k)} \right|}{\tau_0^{(k)}} \right)^{n^{(k)}} \operatorname{sgn}(\tau^{(k)}), \quad \tau^{(k)} = \boldsymbol{\sigma} : \boldsymbol{\mu}^{(k)}, \tag{4}$$

where  $\tau_0^{(k)}$ , the reference resolved shear stress on system k, depends on the activity of the other systems through: 302

$$\dot{\tau}_0^{(k)} = \sum_{\ell=1}^M H^{(k,\ell)} \left( \frac{\tau_{sta}^{(\ell)} - \tau_0^{(\ell)}}{\tau_{sta}^{(\ell)} - \tau_{ini}^{(\ell)}} \right) \left| \dot{\gamma}^{(\ell)} \right|. \tag{5}$$

The two material parameters  $au_{ini}^{(\ell)}$  and  $au_{sta}^{(\ell)}$  refer, respectively, to the initial value of  $\tau_0^{(\ell)}$  at the onset of plasticity (when the  $\gamma^{(k)}$ 's are small ) and to the stationary value of  $\tau_0^{(\ell)}$  at saturation when the plasticity is fully developed 305 (i.e. when the  $\gamma^{(k)}$ 's are large). Therefore the contribution of system  $\ell$  in 306 the hardening (or softening) of system k vanishes when  $\tau_0^{(\ell)}$  is close to  $\tau_{sta}^{(\ell)}$ . 307 The hardening matrix  $H^{(k,\ell)}$  expresses the influence of the plastic activity 308 of system  $\ell$  on the hardening of system k and is taken to be symmetric. 309 Material data for this model are given in Castelnau et al. (2008b). In (Suguet et al., 2011), Eqs (4) and (5) are improved in two ways:

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1. Kinematic hardening is introduced in (4) through a back stress  $X^{(k)}$ :

$$\dot{\gamma}^{(k)} = \dot{\gamma}_0^{(k)} \left( \frac{\left| \tau^{(k)} - X^{(k)} \right|}{\tau_0^{(k)}} \right)^{n^{(k)}} \operatorname{sgn} \left( \tau^{(k)} - X^{(k)} \right), \tag{6}$$

where the back stress evolves with the plastic activity according to an Armstrong-Frederick type law (Chaboche, 2008):

$$\dot{X}^{(k)} = c^{(k)}\dot{\gamma}^{(k)} - d^{(k)}X^{(k)} \left| \dot{\gamma}^{(k)} \right| - e^{(k)}X^{(k)}, \tag{7}$$

including static recovery through coefficient  $e^{(k)}$ . The introduction of a back stress on each slip system is motivated by the experimental observation of recovery strain developing in single crystals when specimens are subjected to recovery tests (see Section 2.3.2 and Fig. 4).

2. The equation governing the reference resolved shear stress  $\tau_0^{(k)}$  is modified into

$$\dot{\tau}_0^{(k)} = \left(\tau_{sta}^{(k)} - \tau_0^{(k)}\right)\dot{p}^{(k)}, \quad \dot{p}^{(k)} = \sum_{\ell=1}^M H^{(k,\ell)} \left|\dot{\gamma}^{(\ell)}\right| . \tag{8}$$

The motivation for the change in the evolution rule for the reference resolved shear stresses  $\tau_0^{(k)}$  is that with the original rule (5) they never reach their stationary value, unless all systems do so at the same time, a condition which cannot be met in a polycrystal (see details in (Suquet et al., 2011)). By contrast, the law (8) ensures convergence of  $\tau_0^{(k)}$  towards its stationary value, provided all coefficients  $H^{(k,\ell)}$  are positive. Indeed, in this case,  $\dot{p}^{(k)}$  is always positive and  $p^{(k)}$  is increasing with time, acting on system k in a similar way as the classical cumulated plastic strain of von Mises plasticity. The differential Eq. (8) can be integrated into

$$\tau_0^{(k)}(p^{(k)}) = \tau_{sta}^{(k)} + (\tau_{ini}^{(k)} - \tau_{sta}^{(k)}) \exp(-p^{(k)}), \tag{9}$$

which shows that  $\tau_0^{(k)} - \tau_{sta}^{(k)}$  has the same sign as  $\tau_{ini}^{(k)} - \tau_{sta}^{(k)}$ . Furthermore  $\tau_0^{(k)}$  tends to  $\tau_{sta}^{(k)}$  when  $p^{(k)}$  becomes large.

## 2.3.1. Data for elasticity

As mentioned in Section 1, ice crystals exhibit a low elastic anisotropy, the largest stiffness ( $E \sim 11.8$ GPa) being along the c-axis (Fig. 3). The tensor of elastic moduli (in Kelvin's notations) at  $-16^{\circ}C$  is given by (10) (Gammon et al., 1983),

$$\begin{pmatrix}
\sigma_{11} \\
\sigma_{22} \\
\sigma_{33} \\
\sqrt{2}\sigma_{23} \\
\sqrt{2}\sigma_{13} \\
\sqrt{2}\sigma_{12}
\end{pmatrix} = \begin{pmatrix}
13930. & 7082. & 5765. & 0. & 0. & 0. \\
7082. & 13930. & 5765. & 0. & 0. & 0. \\
5765. & 5765. & 15010. & 0. & 0. & 0. \\
0. & 0. & 0. & 6028. & 0. & 0. \\
0. & 0. & 0. & 0. & 6028. & 0. \\
0. & 0. & 0. & 0. & 6028. & 0. \\
0. & 0. & 0. & 0. & 6848.
\end{pmatrix}
\begin{pmatrix}
\varepsilon_{11}^{e} \\
\varepsilon_{22}^{e} \\
\varepsilon_{33}^{e} \\
\sqrt{2}\varepsilon_{13}^{e} \\
\sqrt{2}\varepsilon_{13}^{e}
\end{pmatrix}, (10)$$

where all entries are in MPa and 3 is the axis of transverse isotropy (c-axis of the hexagonal crystalline structure). For conditions prevailing in ice sheets and glaciers, elastic constants vary little with temperature: a temperature change of 5°C only modifies the elastic constants by about 1.5%.

## 343 2.3.2. Data for basal slip

The literature provides a number of experimental data for the behavior 344 of ice crystals deformed in such a way that only basal slip is activated. Due 345 to the very large viscoplastic anisotropy of ice single crystals, it is stressed 346 that mechanical tests have to be carried out very carefully to avoid any heterogeneity of the stress field within the specimen (Boehler et al., 1987). 348 Mechanical tests on single crystals where solely non-basal systems are 349 activated have not been reported so far. This would require straining the 350 crystal along or perpendicular to the c-axis, but unfortunately any unavoid-351 able deviation from perfect alignment activates basal slip. Duval et al. (1983) 352 has given upper bounds for the flow stress on non-basal systems. 353

Consequently, only the material parameters of Eq. (6) relevant for basal slip can be identified with confidence from experimental data on single crystals:

- First, data compiled by Duval et al. (1983) were used to determine the stationary flow stress and the stress-sensitivity exponent  $n^{(k)}$  of basal slip. There is quite a large spread in these experimental results from different authors. Despite these uncertainties, the stress-sensitivity exponent for basal slip can be directly identified from these experimental data (numerical values are reported in Table 1), whereas the stationary flow stress depends on both the stationary reference stress  $\tau_{sta}^{(k)}$  and the stationary backstress  $X^{(k)}$ .
- Next, data from Weertman (1973) were used for the identification of the transient creep regime of basal systems. Single crystals were deformed under uniaxial compression at different strain rates, with c-axis oriented at 45° from the loading direction (Fig. 4). The observed stress peak is associated with the increase in density of mobile dislocations (Duval et al., 1983), a behavior typical for material with very low initial dislocation density (see Sauter and Leclercq (2003); Cochard et al. (2010)). These tests shed light on the softening of basal slip in the transient regime. The static recovery term e<sup>(k)</sup> in the constitutive law (7) helps achieving the correct stationary stress at very small strain rates (since X<sup>(k)</sup> tends to a constant value c<sup>(k)</sup>/d<sup>(k)</sup> at large shear γ<sup>(k)</sup> if static recovery is not introduced).
  - Finally, the recovery test of Taupin et al. (2008) performed on single crystals under uniaxial compression was considered. Here the caxis orientation was not precisely defined experimentally, but it made

an angle "less than  $10^{\circ}$ " with the compression direction. The single crystal was submitted to four creep loadings for 30 minutes separated by unloading stages for respectively 1 minute, 10 minutes, and 100 minutes (Fig. 4). Upon reloading, the strain rate is larger than just before the last unloading, indicating that dislocations are rearranging during the time intervals where the specimen is unloaded. This is accounted for in the model by the back stress  $X^{(k)}$ , and by  $e^{(k)}$ .

Fig. 4 shows the good match between the model (constitutive Eq. (6)) with the set of parameters given in Table 1 and these experimental results.

# 390 3. Mean field approaches for the mechanical response of ice poly391 crystals

#### 3.1. Microstructure characterization

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From the mechanical point of view, polycrystalline materials have to be 393 considered as a specific class of composites. They are composed of many 394 grains, with grain size in the range of mm to cm for natural ice. Grains are 395 assembled in a random way, i.e. their size, shape, and lattice orientation do 396 generally not depend on the size, shape, and orientation of the surrounding 397 grains (Fig. 5). Therefore, the microstructure of ice polycrystals can hardly 398 be described exactly in 3-D, unless one makes use of tomography techniques 399 (Rolland du Roscoat et al., 2011). From (2-D) thin sections, one can at best 400 access a statistical characterization of the 3-D grain arrangement e.q. with 401 the help of cross-correlation functions, although the description is generally 402 limited to a few parameters, such as the average grain size and grain shape 403 (aspect ratio). In the Euler orientation space, microstructure description is

based on the distribution of crystal lattice orientations (Orientation Distri-405 bution Function, ODF, or crystallographic texture, often denoted "fabric" 406 in the geophysical community). The complex behavior of polycrystalline 407 materials comes from the anisotropic behavior at the grain scale, closely re-408 lated to the symmetry of the crystal lattice. This is true for all quantities of 409 interest here, such as elasticity, viscoplasticity and thermal dilation. Grains 410 with different lattice orientations react differently to a given stress level. 411 As far as grain boundaries maintain the cohesion of the material, the local 412 stress (i.e. inside a grain) differs from the overall one (the applied stress), 413 leading to a heterogeneous distribution of stress and strain fields within the 414 polycrystal. 415

Most research efforts in the past years have focussed on the understand-416 ing of the build-up of these heterogeneities, in relation with the microstruc-417 ture and local (grain) behavior, since they greatly influence the overall be-418 havior (for ice, see Grennerat et al. (2012) for instance). For instance, plas-419 ticity in a polycrystal can start far below the macroscopic yield stress, as it 420 is sufficient that the local stress reaches the local yield stress somewhere in 421 the structure where stress concentration is large enough, such as along grain 422 boundaries (Brenner et al., 2009). 423

There are basically two strategies to get the mechanical response: meanfield (this section) and full-field (next section) approaches. For both of
them, the key issue is the estimation of the stress or strain localization
(or heterogeneities), in relation to the microstructure and local behavior of
grains. Basically, the problem to be solved is to find an equilibrated stress
field, related to a compatible strain field with the local constitutive relation,
both fields fulfilling the applied boundary conditions. In the following, we
review (not in an exhaustive way) some homogenization techniques used for

the investigation of the mechanical behavior of ice polycrystals.

3.2. Linear thermo-elasticity

For reasons that will become evident below, let us consider the case of thermo-elastic ice polycrystals. The local constitutive relation at point  $(\mathbf{x})$  reads

$$\varepsilon(\mathbf{x}) = \mathbf{S}(\mathbf{x}) : \sigma(\mathbf{x}) + \varepsilon_0(\mathbf{x}),$$
 (11)

with  $\varepsilon_0$  a stress-free thermal strain (e.g. a dilation), due to temperature changes. The local stress  $\sigma(\mathbf{x})$  can be related to the overall stress (applied at the polycrystal scale) by means of the stress-concentration tensor  $\mathbf{B}(\mathbf{x})$ for the purely elastic problem

$$\sigma(\mathbf{x}) = \mathbf{B}(\mathbf{x}) : \bar{\sigma} + \sigma_{\text{res}}(\mathbf{x}),$$
 (12)

with  $\sigma_{\rm res}$  the residual stress, *i.e.* the stress field remaining locally when the overall load is suppressed ( $\bar{\sigma} = 0$ ). It can be shown that the overall polycrystal behavior takes a similar form as Eq. (11)

$$\bar{\boldsymbol{\varepsilon}} = \tilde{\mathbf{S}} : \bar{\boldsymbol{\sigma}} + \tilde{\boldsymbol{\varepsilon}}_0 , \qquad (13)$$

with symbols  $\tilde{\cdot}$  and  $\bar{\cdot}$  denoting the homogenized (or effective) property and the volume average over the whole polycrystal volume (also denoted  $\langle . \rangle$ ), respectively. Therefore, one has  $\bar{\sigma} = \langle \sigma(\mathbf{x}) \rangle$  and  $\bar{\varepsilon} = \langle \varepsilon(\mathbf{x}) \rangle$ , and it can be shown that the effective compliance  $\tilde{\mathbf{S}}$  and the effective thermal strain  $\tilde{\varepsilon}_0$ are given by (Laws, 1973)

$$\tilde{\mathbf{S}} = \langle \mathbf{S}(\mathbf{x}) : \mathbf{B}(\mathbf{x}) \rangle, \quad \tilde{\varepsilon}_0 = \langle \varepsilon_0(\mathbf{x}) : \mathbf{B}(\mathbf{x}) \rangle.$$
 (14)

Since, for thermo-elastic polycrystals, the elastic compliance and the thermal dilation coefficients are uniform properties inside grains, the quantities  $\mathbf{S}(\mathbf{x})$  and  $\boldsymbol{\varepsilon}_0(\mathbf{x})$  in Eq. (11) can be replaced by the corresponding

homogeneous values  $\mathbf{S}^{(r)}$  and  $\boldsymbol{\varepsilon}_0^{(r)}$  of the considered phase (r). A similar substitution can be made in Eq. (14), leading to 453

$$\tilde{\mathbf{S}} = \sum_{r} c^{(r)} \mathbf{S}^{(r)} : \bar{\mathbf{B}}^{(r)} , \quad \tilde{\boldsymbol{\varepsilon}}_{0} = \sum_{r} c^{(r)} \boldsymbol{\varepsilon}_{0}^{(r)} : \bar{\mathbf{B}}^{(r)}$$
 (15)

with  $\bar{r}^{(r)}$  indicating the average over the volume of phase (r), e.g.  $\bar{\mathbf{B}}^{(r)} =$  $\langle \mathbf{B}(\mathbf{x}) \rangle^{(r)}$ , and  $c^{(r)}$  the volume fraction of phase (r). Here, a mechanical 455 phase (r) denotes the set of all grains of the polycrystal having the same 456 crystal orientation; those grains have different shape and environment but 457 their elastic and thermal properties are identical. From (15), it can be 458 observed that the sole knowledge of the mean (phase average) values  $\bar{\mathbf{B}}^{(r)}$  is 459 sufficient to estimate the overall polycrystal behavior. It can be anticipated 460 that, if the quantities  $\bar{\mathbf{B}}^{(r)}$  can be calculated without having to know the 461 complete field of  $\mathbf{B}(\mathbf{x})$ , computation will be way faster. Hence the name of 462 "mean-field" approaches presented here. 463 With the effective behavior (Eq. 14) in hand, statistical averages over 464 crystal orientations (r) can be estimated. Basically, two quantities can be 465 obtained from mean-field approaches: 466

1. The phase average stress (or first moment)  $\bar{\sigma}^{(r)} = \langle \sigma(\mathbf{x}) \rangle^{(r)}$ 

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$$\bar{\boldsymbol{\sigma}}^{(r)} = \bar{\mathbf{B}}^{(r)} : \bar{\boldsymbol{\sigma}} + \bar{\boldsymbol{\sigma}}_{res}^{(r)}, \tag{16}$$

with  $\bar{\sigma}_{res}^{(r)}$  the average residual stress of phase (r). The knowledge 468 of  $\bar{\sigma}^{(r)}$  for all phases (r) allows investigating the so-called *interphase* heterogeneities, i.e. the variation of the phase average stress with 470 respect to the crystal orientation. 471

2. Deeper insight into the stress distribution can be obtained from the second moment  $\langle \boldsymbol{\sigma} \otimes \boldsymbol{\sigma} \rangle^{(r)}$  of the stress. This second moment can be obtained by a derivation of the effective energy with respect to local compliances, see (Bobeth and Diener, 1987; Kreher, 1990; Ponte-Castañeda and Suquet, 1998; Brenner et al., 2004).

The standard deviation of the stress distribution within a given crystal ori-477 entation (r) can be estimated from these two moments as the square root 478 of  $<\sigma\otimes\sigma>^{(r)}-<\sigma>^{(r)}\otimes<\sigma>^{(r)}$  (i.e. the mean of the square of 479 the stress minus the square of the mean); it is related to the width of the 480 stress distribution in crystal orientation (r), and accounts for both the het-481 erogeneity of stress distribution *inside* grains but also for the heterogeneity 482 between grains of identical orientation but exhibiting different shapes and 483 having different neighborhood. Similar relations can be derived for the strain statistics. 485

#### 486 3.3. Reuss and Voigt approximations

First, two very basic models can be derived, namely Reuss (also called 487 static in the viscoplastic context) and Voigt (or Taylor) models. The Reuss model is constructed by considering a uniform stress throughout the poly-489 crystal, i.e.  $\sigma(\mathbf{x}) = \bar{\sigma} \ \forall \mathbf{x}$ , or equivalently  $\mathbf{B}(\mathbf{x}) = \mathbf{I}$  (with I the identity ten-490 sor), and leads to vanishing intra- and inter-granular stress heterogeneities, 491 and uniform strain within grains. The Voigt model considers uniform strain, 492 i.e.  $\varepsilon(\mathbf{x}) = \bar{\varepsilon} \ \forall \mathbf{x}$ , i.e. no intra– and inter–granular strain heterogeneities, 493 and uniform stress within grains. These models violate strain compatibility 494 and stress equilibrium, respectively, and are of limited accuracy when the 495 local behavior is highly nonlinear and/or highly anisotropic, as will be il-496 lustrated in the next section. Besides simplicity, the main interest of Reuss 497 and Voigt models is based on their bounding character, since they provide, 498 respectively, a lower and an upper bound for the effective stress potential. 499

## 3.4. The Self-Consistent (SC) scheme

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Unlike full-field approaches detailed in the Section 4, mean-field methods 501 are based on a statistical description of the microstructure, e.g. based on 502 few n-points correlation functions, so that the exact position and shape of 503 a specific grain with respect to its neighbors is not known. However, as al-504 ready introduced, all grains exhibiting the same crystallographic orientation 505 are treated as a single mechanical phase. Owing to the random character 506 of the microstructure with all grains playing geometrically similar roles, the 507 Self-Consistent (SC) scheme (Hershey, 1954; Kröner, 1958; Willis, 1981) is 508 especially well suited for polycrystals. This model, which provides a rela-509 tively simple expression for  $\bar{\mathbf{B}}^{(r)}$ , relies on specific microstructures exhibiting 510 perfect disorder and infinite size graduation (Kröner, 1978). The SC scheme 511 has often been described as if the interaction between each grain and its surrounding could be approximated by the interaction between one ellipsoidal 513 grain with the same lattice orientation as the original grain and a homoge-514 neous equivalent medium whose behavior represents that of the polycrystal, taking thus advantage of the analytical solution of Eshelby (1957) for the in-516 clusion/matrix interaction. This reasoning led to the conclusion that the SC 517 scheme implicitly considers uniform stress and strain rate inside grains. This 518 interpretation turns out to be incorrect, since intraphase stress and strain 519 heterogeneities do not vanish as explained above, see Ponte-Castañeda and 520 Suguet (1998) for a review. 521 The ability of the SC scheme to estimate polycrystal behavior is shown 522 in Fig. 6. Numerical reference solutions from the full-field FFT method 523 (see Section 4) have been generated for many randomly generated Voronoi 524 microstructures, and ensemble average over these random microstructures

has been calculated in order to attain results that are representative for

a Representative Volume Element, *i.e.* a volume sufficiently large to be statistically representative of the material (Kanit et al., 2003; Lebensohn et al., 2004b). In Fig. 6, we provide results for the effective behavior, that is entirely defined by the effective reference stress  $\tilde{\sigma}_0$  which enters in the effective constitutive relation

$$\frac{\dot{\bar{\varepsilon}}_{eq}}{\dot{\varepsilon}_0} = \frac{\bar{\sigma}'_{eq}}{\tilde{\sigma}_0} \tag{17}$$

with  $\dot{\varepsilon}_0$  a reference strain rate (taken here equal to  $\dot{\gamma}_0$ ), and  $\bar{\sigma}'_{eq}$  and  $\dot{\bar{\varepsilon}}_{eq}$  the effective equivalent stress and strain rate respectively  $(\bar{\sigma}'_{eq} = \sqrt{3\bar{\sigma}: \bar{\sigma}/2}, \dot{\bar{\varepsilon}}_{eq} =$ 533  $\sqrt{2\dot{\bar{\varepsilon}}:\dot{\bar{\varepsilon}}/3}$ ). Calculations are performed for various viscoplastic anisotropy 534 contrasts (or slip system contrasts) at the grain level, defined by the ratio 535 between non-basal and basal reference shear stresses, i.e.  $au_0^{(Pr)}/ au_0^{(Ba)}$  for 536  $\tau_0^{(Pyr)} = \tau_0^{(Pr)}$ . It can be observed that the SC model perfectly reproduces the reference full-field (FFT) results. Note also that the Reuss bound, of-538 ten used for highly anisotropic materials like ice, predicts a much too soft 539 overall behavior. This simple approach does not allow to make a realistic link between local and overall rheologies. We also report in this figure 541 the standard deviations (or overall heterogeneities) of equivalent stress and 542 strain rate. These standard deviations have been calculated over the whole 543 polycrystal. Recall that they account for both intra- and inter-granular field 544 heterogeneities for both SC and FFT approaches. It can be observed that 545 the increase of standard deviation with the slip system contrast is well re-546 produced by the SC scheme, although some discrepancies with FFT results arise at very large contrasts (mostly for the strain-rate fluctuation). Note 548 again that Reuss and Voigt bound do not reproduce these results, even in 549 a qualitative way, since they predict, by construction, vanishing fluctuation of stress and strain rate, respectively. Unlike these simple approaches, the

SC scheme not only predicts the correct effective stress, but also accurately captures the field heterogeneities within the polycrystal. Similar agreement have been obtained for Voronoi and EBSD 2-D microstructures under antiplane shear by Lebensohn et al. (2005).

#### 556 3.5. Nonlinear viscoplasticity

The mean-field estimate of *nonlinear materials* is significantly more complex than the thermo-elastic case treated above. We consider the case of a viscoplastic polycrystal of ice in which grains are deforming by glide of dislocations on specific slip planes, as discussed above, with slip rates given by Eq. (4), so that the local strain rate reads, since elastic deformations are neglected:

$$\dot{\varepsilon}(\mathbf{x}) = \sum_{k} \mu_{(k)}^{(r)} \dot{\gamma}_{(k)}(\mathbf{x}) . \tag{18}$$

Here, reference stresses  $\tau_0$  and stress sensitivities n are supposed to be constant. The constitutive Eq. (18) can also be written

$$\dot{\boldsymbol{\varepsilon}}(\mathbf{x}) = \mathbf{M}(\mathbf{x}) : \boldsymbol{\sigma}(\mathbf{x}) \tag{19}$$

565 with

$$\mathbf{M}(\mathbf{x}) = \sum_{k} \frac{\dot{\gamma}_0^{(k)}}{\tau_0^{(k)}} \left| \frac{\boldsymbol{\mu}^{(k)}(\mathbf{x}) : \boldsymbol{\sigma}(\mathbf{x})}{\tau_0^{(k)}} \right|^{n^{(k)} - 1} \boldsymbol{\mu}^{(k)}(\mathbf{x}) \otimes \boldsymbol{\mu}^{(k)}(\mathbf{x}) . \tag{20}$$

Obviously, the viscous compliance  $\mathbf{M}$  relating  $\dot{\boldsymbol{\varepsilon}}(\mathbf{x})$  and  $\boldsymbol{\sigma}(\mathbf{x})$  – which plays a similar role as  $\mathbf{S}$  in Eq. (11) – is not uniform within a phase, owing to the stress sensitivities  $n \neq 1$  and the heterogeneity of  $\boldsymbol{\sigma}$  in the phases. Consequently, (14) cannot be replaced by (15) for nonlinear behavior. The basic method to deal with such nonlinear behavior is to define a *Linear Compari*son *Polycrystal* (LCP) having the same microstructure as the real nonlinear

polycrystal, and to which the linear homogenization scheme applies (Ponte-572 Castañeda and Suguet, 1998). Of course, the effective behavior estimated 573 this way remains nonlinear, since the definition of the LCP depends on the 574 applied macroscopic stress. The difficult part of the problem consists of 575 finding the best *linearization* procedure leading to the optimal selection of 576 the LCP. Since decades, there has been quite a number of propositions in the literature dealing with this issue, leading to a generalization of the SC 578 scheme for nonlinear behavior. The local constitutive relation given by Eqs. 579 (18-20) has to be linearized in a suitable way to obtain a form similar to 580 (11), with **S** and  $\dot{\varepsilon}_0$  uniform per phase (and where  $\varepsilon$  is replaced everywhere 581 by  $\dot{\varepsilon}$ ). Generally speaking, the linearization can be expressed in the form 582 depicted in Fig. 7, (Liu and Ponte Castañeda, 2004).

$$\dot{\gamma}_{(k)}(\mathbf{x}) = \alpha_{(k)}^{(r)} \tau_{(k)}(\mathbf{x}) + \dot{e}_{(k)}^{(r)}, \tag{21}$$

thus leading to the following expressions for  $\mathbf{S}^{(r)}$  and  $\dot{oldsymbol{arepsilon}}_0^{(r)}$ 

$$\mathbf{S}^{(r)} = \sum_{k} \alpha_{(k)}^{(r)} \boldsymbol{\mu}_{(k)}^{(r)} \otimes \boldsymbol{\mu}_{(k)}^{(r)}, \qquad \dot{\boldsymbol{\varepsilon}}_{0}^{(r)} = \sum_{k} \dot{\boldsymbol{e}}_{(k)}^{(r)} \boldsymbol{\mu}_{(k)}^{(r)}, \qquad (22)$$

where the shear compliance  $\alpha_{(k)}^{(r)}$  and stress-free slip-rate  $\dot{e}_{(k)}^{(r)}$  can be easily expressed with respect to two reference shear stresses  $\check{\tau}_{(k)}^{(r)}$  and  $\hat{\tau}_{(k)}^{(r)}$ , see Fig. 7. The optimal choice (from the point of view of the variational mechanical problem) of those reference stresses is not straightforward; this is the main reason why several extensions of the SC scheme have been proposed in the literature. Obviously, all of them reduce to the same SC model in the linear case n=1.

Following Ponte Castañeda (1996), Masson et al. (2000) proposed the

so-called "affine" (AFF) linearization scheme which is based on the simple

idea of a linear behavior (21) tangent to the nonlinear one (4) at the mean shear stress, leading to

The main limitations of this procedure are discussed in detail in Masson

$$\check{\tau}_{(k)}^{(r)} = \hat{\tau}_{(k)}^{(r)} = \langle \tau_{(k)} \rangle^{(r)}, \quad \alpha_{(k)}^{(r)} = \left. \frac{\partial \dot{\gamma}}{\partial \tau} \right|_{\tau = \check{\tau}_{(k)}^{(r)}}.$$
(23)

et al. (2000) and Bornert and Ponte Castañeda (1998). One of them is the 597 violation of rigorous upper bounds for the effective behavior. More generally, 598 the affine extension is known to overestimate the overall viscosity, i.e. to 590 predict an effective behavior that is too stiff. This negative feature can be 600 alleviated by means of the energy formulation originally proposed by Ponte 601 Castañeda (1996) (see Bornert et al. (2001)). 602 603 Alternative, more sophisticated ways to generalize the SC scheme have been proposed by Ponte Castañeda and co-workers during the last decades. 604 The basic idea of these methods is to guide the choice of the properties of the 605 LCP by a suitably designed variational principle. An "optimal" solution has 606 been obtained in the context of the so-called "variational" procedure (VAR) 607 (Ponte Castañeda, 1991), which was extended to polycrystals by De Botton 608 and Ponte Castañeda (1995), leading to the choice 609

$$\check{\tau}_{(k)}^{(r)} = 0, \quad \hat{\tau}_{(k)}^{(r)} = \left[ \langle \tau_{(k)}^2 \rangle^{(r)} \right]^{1/2} .$$
(24)

The main advantage of this procedure is to provide a rigorous bound, sharper than the Voigt bound, for the effective potential. More recently, the "second-order" (SO) method of Ponte Castañeda (2002), extended to polycrystals in (Liu and Ponte Castañeda, 2004) has been proposed. It leads to reference shear stresses reading

$$\check{\tau}_{(k)}^{(r)} = \langle \tau_{(k)} \rangle^{(r)}, \quad \hat{\tau}_{(k)}^{(r)} = \check{\tau}_{(k)}^{(r)} \pm \left[ \langle (\tau_{(k)} - \check{\tau}_{(k)}^{(r)})^2 \rangle^{(r)} \right]^{1/2}. \tag{25}$$

The main differences between AFF, VAR, and SO models may be sum-615 marized as follows. The AFF estimate can be regarded as a relatively simple 616 model, allowing rapid computations which can even be rather accurate for 617 polycrystals with weak grain anisotropy and small stress sensitivity. How-618 ever, its predictions can become unrealistic (e.g. bound violation) at a strong 619 anisotropy or nonlinearity. Contrary to AFF, for which linearization only accounts for the phase average stress, VAR accounts for the second moments 621 of the stress, whereas the SO procedure accounts for both the phase aver-622 age stress and intraphase standard deviation (first and second moments) 623 to build the LCP. They can therefore provide better estimates in cases of 624 highly heterogeneous stress distributions, such as for strongly nonlinear or 625 anisotropic polycrystals. Applications of the VAR procedure to polycrystals 626 with grains having cubic or hexagonal crystallographic structures can be 627 found in (Nebozhyn et al., 2001; Liu et al., 2003). 628

Finally, the "tangent" (TGT) extension of the SC scheme (Molinari 629 et al., 1987; Lebensohn and Tomé, 1993), often referred to as the "VPSC 630 model" in the literature, is based on the same tangent linearization (23) as 631 the AFF method. However, unlike the AFF extension, this procedure takes 632 advantage of the fact that, for power-law polycrystals with a single stress 633 exponent n for all slip systems, the tangent behavior (21) can be replaced 634 by a secant-like relation, with  $\dot{e}_{(k)}^{(r)}=0$  and  $\alpha_{(k)}^{(r)}$  replaced by  $\alpha_{(k)}^{(r)}/n$ . The 635 same procedure is further applied at the macroscopic level, leading to an 636 inconsistent definition for the stress localization tensor  $\mathbf{B}^{(r)}$  that combines a 637 secant description for the local and global behaviors but a tangent analysis 638 for the inclusion/matrix interaction (Masson et al., 2000). When expressed 639 in the form of tangent expressions, it can be shown that  $\dot{\tilde{\varepsilon}}_0$  differs from the exact relation given in (15).

#### 3.5.1. Application to natural ices: effective behavior

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Application of homogenization techniques to natural ices aims at un-643 derstanding (and predicting) the anisotropic behavior of strongly textured 644 specimens, as encountered at depth in natural ice sheets. As will be seen in 645 section 6, the viscoplastic anisotropy of polycrystals significantly influences ice flow at large scales (Mangeney et al., 1996; Gillet-Chaulet et al., 2006; 647 Pettit et al., 2007; Martín et al., 2009). Castelnau et al. (1998) reported mechanical tests performed on specimens from the GRIP ice core (Central 649 Greenland). Along the ice core, the ice microstructure, and in particular 650 the crystallographic fabric, is evolving; with increasing depth, randomly ori-651 ented c-axis at the surface of the ice sheet tend to concentrate towards the 652 in situ vertical direction down to a depth of  $\sim 2600$  m. Beneath this depth, 653 less pronounced textures are observed due to the initiation of migration re-654 crystallization (Thorsteinsson et al., 1997). In (Castelnau et al., 1998), the experimental stationary creep behavior of those ices have been obtained for 656 two loading conditions (Fig. 8). The first one corresponds to an in situ ver-657 tical compression, showing an increasing flow stress (decreasing strain rate for a constant applied stress) with increasing depth, since the activation of 659 non-basal slip systems is necessary for pronounced fabrics. The second load-660 ing condition corresponds to in situ horizontal shear, promoting basal slip 661 and resulting in a softening of the ice with increasing fabric strength. It can 662 be seen that for a given applied stress, strain rates can vary by more than 663 two orders of magnitude depending on the orientation of the applied stress 664 with respect to the specimen fabric, reflecting the very strong viscoplastic 665 anisotropic of ice specimens. 666

The effective behavior predicted by the affine (AFF) SC model is com-

pared to the experimental data in Fig. 8. It can be observed that the agree-668 ment is excellent, meaning that the relation between fabric and effective rhe-669 ology is very well captured by the model. The model captures correctly the 670 increasing anisotropy from the surface down to  $\sim 2600$  m depth, and the de-671 crease below. The difference by more than two orders of magnitude between 672 the vertical and shear strain-rates at  $\sim 2600$  m is also well reproduced, although this was a challenging feature for the model. To get these results, the 674 reference shear stress  $\tau_0^{(k)}$  entering the local constitutive relation, and also 675 the stress sensitivity  $n^{(k)}$ , for each slip system (k), had to be identified from 676 comparison with a database that included single-crystal experimental tests, 677 and polycrystal ones on many different crystallographic textures (Castelnau 678 et al., 2008b). The resulting single-crystal rheology, used as input in the SC 679 model to get the effective behavior described above, is shown in Fig. 9. For basal slip, agreement with experimental data from the literature is almost 681 perfect. Non-basal systems are much stiffer than the basal systems, and 682 pyramidal slip is found to be much more difficult than prismatic slip. These results are in good agreement with the available data on single crystals, and in qualitative agreement with the known dislocation structure in ice. There-685 fore, it can be anticipated that the affine SC model does a good job in making 686 the link between the grain and the polycrystal scales, and provides an ac-687 curate estimate of the mechanical interaction between deforming grains. In 688 other words, one can anticipate that results shown in Fig. 8 are based on a 689 realistic description of the mechanical interaction between grains and physical deformation processes (dislocation glide) at the (sub)grain level. It can 691 also be seen on Fig. 9 that this identification procedure leads to different 692 stress sensitivities for the different slip system families. A value  $n^{(k)}=2$  was imposed for basal slip in accordance with experimental data, but values for

prismatic and pyramidal systems were considered as adjustable parameters. It is also worth noting that the affine model perfectly reproduces an effective 696 stress sensitivity (i.e. at the polycrystal scale)  $\tilde{n}=3$  in agreement with experimental data, although the two major slip systems, basal and prismatic 698 slip, have stress sensitivities smaller than 3  $(n^{(bas)} = 2.0, n^{(pr)} = 2.85)$ . A 699 larger value was considered only for pyramidal slip  $(n^{(py)} = 4.0)$ , but it is 700 worth mentioning that the contribution of pyramidal slip is only very mi-701 nor ( $\langle 2\% \rangle$ ). It can be concluded that, in ice, although basal slip is by far 702 the most active deformation mechanism, secondary slip systems are of great 703 importance for explaining the polycrystal behavior. Basal slip alone does 704 not allow for plastic deformation of ice polycrystals, since it only provides 705 two independent slip systems. Secondary slip systems, here prismatic and 706 pyramidal slip, must therefore be activated to add two more independent 707 slip systems. The strength of these stiffer mechanisms determines the vis-708 coplastic anisotropy at the grain scale, and therefore they also control the 709 level of inter—and intra—granular heterogeneities of stress and strain(-rate), and therefore the effective polycrystal rheology. Similar conclusions have 711 been drawn for olivine, a mineral with only three independent slip systems 712 (Castelnau et al., 2008a, 2009, 2010a,b). We therefore anticipate that the 713 strong effect of secondary deformation mechanisms observed here might be 714 a general feature for all polycrystalline materials with less than four inde-715 pendent slip systems. The corollary of these results is that simple or ad hoc 716 polycrystal models, such as the Reuss (uniform stress) model, in which ice polycrystals can deform with only basal slip, cannot be accurate. This has 718 been shown for example in (Castelnau et al., 1997): whatever the strength 719 used for prismatic and pyramidal systems, the Reuss model is not able to reproduce the very large anisotropy of GRIP specimens shown in Fig. 8.

This comes from the fact that internal stresses, that have a large influence on the material behavior, are ignored.

Finally, it is also worth mentioning that the TGT SC approach, used 724 in earlier studies, e.g. (Castelnau et al., 1997), does not provide as good a 725 match to experimental data as the AFF SC extension. There can be two 726 reasons for that: (i) first of all, it is now known that the inconsistency in 727 the formulation of the TGT SC version leads to an underestimation of the 728 internal stress level, predicting a too soft polycrystal behavior (Gilormini, 729 1995; Masson et al., 2000); (ii) second, by construction, the TGT model is 730 limited to grain behavior for which all slip systems exhibit the same stress 731 sensitivity  $n^{(k)}$ . When applied to ice, one must thus consider  $n^{(k)} = 3.0$  for 732 all systems, including basal slip, in order to get an effective  $\tilde{n}=3$ . The fact 733 that the AFF extension does not have this limitation might also explain a better consistency with experimental data. 735

#### 736 3.5.2. Application to natural ices: texture development

Using the Reuss approximation, Van der Veen and Whillans (1994) and 737 Castelnau and Duval (1994) described the fabric evolution under compres-738 sion, tension, simple and pure shear. Van der Veen and Whillans (1994) 739 needed to impose a kind of "recrystallization" criterion (see Section 5.1) to be able to correctly represent the single-maximum fabric (with c-axis ori-741 ented along one direction) in ice deforming in pure shear. Nevertheless, the 742 Reuss approximation faces inconsistency to describe the fabric evolution at the polycrystal scale, as it requires additional kinematical constraints to 744 link the grain rotation-rate with the polycrystal rotation-rate. In most of 745 the "Reuss" type models, these two rates are supposed to be equal, although 746 the velocity field is not continuous.

Models that modify this homogeneous stress assumption were proposed 748 by Azuma (1994) and Thorsteinsson (2002). They introduce some redistri-749 bution of stress through neighborhood interaction to define the crystal strain 750 at a given bulk equivalent strain. In particular, Thorsteinsson (2002) defines 751 a crystal arrangement on a three-dimensional cubic grid, where each crystal 752 has six nearest neighbors. The nearest neighbor interaction (NNI) is taken 753 into account by defining a local softness parameter for each crystal which 754 modifies the stress acting on the central crystal compared to the macro-755 scopic stress. This softness parameter further influences the rotation rate 756 of the crystal lattice compared to the bulk. For uniaxial compression tests, 757 the fabrics obtained with the NNI formulation are less concentrated than 758 the ones where no NNI is considered. The reason for this is that the NNI 759 formulation allows all crystals to deform to some extent, while only "soft" 760 crystals would deform in the no-NNI formulation. The fabric obtained after 761 50% shortening strain compares qualitatively well with the one measured 762 along the GRIP ice core at a depth where the strain is similar (1293 m) 763 (Thorsteinsson et al., 1997). 764

The VPSC model in its "tangent" version was applied to simulate the 765 fabric development along ice cores (Castelnau et al., 1996b,a, 1998). In 766 (Castelnau et al., 1996a), a comparison was made with bound estimates 767 (Reuss and Voigt). Fabrics simulated in uniaxial compression and extension 768 were found to be qualitatively similar for all models. However, large dif-769 ferences in the rate of fabric development were found. This was explained 770 by the different interaction stiffness between grain and matrix for the three 771 approaches. The fabrics obtained with the VPSC model for uniaxial defor-772 mation were in close agreement with the one measured along the ice core (see Fig. 10). In particular, this model well reproduced the fabric evolution

along the GRIP ice core within the upper 650 m where dynamic recrystallization is not supposed to strongly impact this evolution (Castelnau et al., 776 1996b). Lower down, the modeled fabric concentration is too high. Although Castelnau et al. (1996b) attributed this discrepancy to the effects 778 of rotation recrystallization along the core, it was later shown that the tan-779 gent approximation overestimates the lattice rotation. In simple shear, the 780 single-maximum fabric found along the ice cores or experimentally could 781 not be reproduced with the VPSC scheme. To get close to this fabric, an 782 extensive (and probably unrealistic) activity of non-basal slip systems was 783 required. More recently, the "second order" (SO) mean field method of 784 Ponte Castañeda (2002) was used to simulate the fabric development along 785 the Talos Dome ice core (Montagnat et al., 2012). Although no recrystalliza-786 tion mechanisms were implemented in this version, the fabric development 787 was astonishingly well reproduced, under the crude assumption of uniaxial 788 compression with a constant strain rate (see Fig. 11). In particular, a good 789 match was obtained when the initial fabric is non isotropic and similar to the one measured in the top firn, at 18 m depth. The cumulated compres-791 sive strain along the core was derived from the thinning function provided 792 by the TALDICE-1 chronology (Buiron et al., 2011). The good prediction 793 performed by the VP-SO scheme is probably due to the fact that this SO ap-794 proach provides a better estimate of the effective behavior than the classical 795 tangent "VPSC" model does in the case of strongly anisotropic materials 796 such as ice (see Section 3.5). Nevertheless, the modeled fabric evolution 797 could not capture the strengthening rate associated with the Glacial to In-798 terglacial climatic transition. At these transition, a change in ice viscosity 799 is expected. It induces an higher sensitivity to the impact of shear stress increasing with depth, that the modeling approach did not considered.

It is also recalled that the heterogeneity of shear on slip systems at the 802 grain level gives rise to heterogeneities of lattice rotation, and therefore 803 generates intragranular misorientations that somehow spread crystal orien-804 tations. It is however worth mentioning that all models presented above 805 do not consider this strain heterogeneity for estimating fabric evolutions at 806 finite strain. Even in VAR and SO procedures, intraphase strain hetero-807 geneities are considered for defining the LCP, but so far not for estimating 808 microstructure evolutions. As a consequence, mean-field approaches gener-809 ally predict too sharp textures. The same applies to the prediction of strain 810 hardening, associated with dislocation processes such as storage and an-811 nealing. A quantitative study, based on comparisons with reference results 812 obtained by a FFT full-field approach, can be found in (Castelnau et al., 813 2006). 814

Most of the efforts to simulate the fabric development in ice, and especially along ice cores, had to face the fact that recrystallization mechanisms could impact this fabric development. This was, most of the time, the analysis made for the observed discrepancies between simulated and measured fabrics (Van der Veen and Whillans, 1994; Wenk et al., 1997; Castelnau et al., 1996b; Thorsteinsson, 2002). Some efforts to implement recrystallization mechanisms in mean-field approaches will be described in Section 5.1.

#### 3.6. Modeling the elasto-viscoplastic behavior

Transient creep is typically encountered when ice flow changes direction, such as in glaciers flowing above irregular bedrock or submitted to tide forcing close to the sea-shore or in icy satellites. During laboratory experiments, transient creep is characterized by a strain-rate drop of more than two orders of magnitude before reaching the secondary creep close to 1% strain,
following Andrade's law (Duval, 1978). This decrease is associated with the
development of large internal stress fields due to intergranular interactions
and a strong kinematic hardening (Duval et al., 1983; Ashby and Duval,
1985; Castelnau et al., 2008b). To reproduce this transient behavior, one
has to consider the coupling between elasticity and viscoplasticity that gives
rise to the so-called "long-term memory effect", as explained below.

The application of homogenization schemes to the *elasto-viscoplasticity* 835 of polycrystals is more complicated than for viscoplasticity, see for instance 836 (Laws and McLaughlin, 1978). In short, it can be shown that, even in the 837 simple case of a polycrystal comprising grains whose behavior exhibits a sin-838 gle relaxation time (so-called "short-term memory"), the effective behavior 839 exhibits a continuous spectrum of relaxation time ("long-term memory ef-840 fect") (Sanchez-Hubert and Sanchez-Palencia, 1978; Suquet, 1987). In other 841 words, the overall behavior of a polycrystal is not of the Maxwell type (par-842 allel association of a spring and a dashpot with constant viscosity), even though the individual grains do exhibit local Maxwell type behavior. The 844 basic difference between elasto-viscoplasticity and viscoplasticity is that, for 845 elasto-viscoplasticity, the local strain rate depends on both the stress (vis-846 cous part) and the stress-rate (elastic part), whereas it only depends on the 847 stress for viscoplasticity. Therefore, the local strain rate not only depends 848 on the actual local stress, but also on the whole stress history from the 849 initial specimen loading at t=0 up to the current time. To obtain the 850 exact effective mechanical response at time t, it is thus required to keep 851 track of all information (or internal variables) corresponding to the strains 852 at all previous times, and therefore the problem is not simple. Within meanfield approaches, some approximations (with hopefully limited effects on the

accuracy of the solution) are thus necessary.

Basically, two approaches have been proposed to deal with this issue. A 856 promising method based on an incremental variational procedure has been proposed by Lahellec and Suquet (2006, 2007). These authors have shown 858 that the homogenization of a linear visco-elastic material (i.e. with n=1) 859 can be expressed in terms of a homogenization problem for a linear ther-860 moelastic composite with non piecewise uniform eigenstrains. One advantage 861 of this formulation is that it can make use of the intraphase heterogeneities 862 of stress and strain (-rate), and it can therefore probably provide accurate 863 results even at high stress sensitivity and/or local anisotropy. An alternative 864 approach, which provides a good compromise between accuracy of the solu-865 tion and simplicity of the formalism, is the so-called "affine" Self-Consistent 866 method of Masson and Zaoui (1999). It is based on the correspondence 867 principle (Mandel, 1966), which states that the elasto-viscoplastic problem 868 can be reduced to a simpler homogenization problem (in fact similar to a 869 standard thermo-elastic problem) if solved in Laplace space. One difficulty of this approach is the calculation of the inverse Laplace transforms, that 871 has to be carried out numerically. An approximate inversion procedure, 872 adapted for creep, has been proposed by Brenner et al. (2002b). It has provided promising results for the creep behavior of Zirconium alloys (Letouzé 874 et al., 2002; Brenner et al., 2002a), since it retains the long-term memory 875 effect associated with the elasto-viscoplastic coupling. Recent developments 876 (Ricaud and Masson, 2009) have shown that an internal variable formulation 877 arises naturally from this affine method, providing results in perfect match 878 with reference FFT solutions in the case of linear viscoelasticity (Vu et al., 879 2012). 880

To the best of our knowledge, this affine method is the only mean-field

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approach that has been applied to simulate the transient creep of ice (Castelnau et al., 2008b). Applications make use of the crystal plasticity model 883 for single crystals detailed in Section 2.3. It was shown that the strong 884 hardening amplitude during the transient creep (i.e. the decrease of the 885 overall strain rate by several orders of magnitude) is explained by the stress 886 redistribution within the specimen: when the overall stress is applied in-887 stantaneously, the instantaneous polycrystal response is purely elastic, and 888 since the elastic anisotropy is small, stress distribution within and between 889 grains is almost uniform. But plastic deformation comes into play rapidly to 890 cause a strong redistribution of stress (with large interphase and intraphase 891 heterogeneities) due to the strong viscoplastic anisotropy at the grain scale. 892 This significantly reduces the overall strain rate. On the other hand, the 893 experimental hardening rate (i.e. the time necessary to reach the secondary 894 creep regime) is much too slow to be explained by the same process, and is 895 attributed to the hardening of hard-glide slip systems (prismatic slip) in the 896 transient regime, associated with dislocation processes (Fig. 12). 897

#### 898 4. Full field approaches for the polycrystal

Mean-field approaches have been extensively used to predict the mechanical behavior of ice polycrystals, and the fabric development as measured
along ice cores. Due to its high viscoplastic anisotropy, deformation in ice is
expected to be strongly heterogeneous, with a strong impact of grain interactions and kinematic hardening (Duval et al., 1983; Hamman et al., 2007;
Montagnat et al., 2011; Grennerat et al., 2012). The mean-field approaches
described above are based on the statistical characterization of the intragranular mechanical fields (in terms of average grain stresses and strain rates,

and, in the most advanced formulations, also through the determination of 907 the intracrystalline average field fluctuations), but the actual micromechan-908 ical fields remain inaccessible to these homogenization approaches. Model-909 ing the full intracrystalline heterogeneity that develops in ice polycrystals 910 requires the use of full-field approaches. This part will concentrate on full-911 field approaches that are using the Fast Fourier Transform method to solve 912 the constitutive equations in a discretized polycrystal. It aims at studying 913 the correlation between the heterogeneous deformation patterns that appear 914 inside the constituent single-crystal grains of an ice aggregate and their cor-915 responding crystallographic orientations, along with the influence of other 916 factors, such as orientation and size of neighboring grains. Both viscoplastic 917 and elasto-viscoplastic behavior were investigated, and are presented in the 918 two following sections. 919

# 920 4.1. Viscoplastic approach - FFT

## 921 4.1.1. Viscoplastic FFT-based formulation

The intracrystalline states that are developed during creep of polycrys-922 talline ice can be obtained using an extension of an iterative method based on 923 FFT, originally proposed by Moulinec and Suquet (1998) and Michel et al. 924 (2001) for linear and non-linear composites (Lebensohn et al., 2009; Mon-925 tagnat et al., 2011). This formulation was later adapted to polycrystals and 926 applied to the prediction of texture development of fcc materials (Leben-927 sohn, 2001), and in turn used for the computation of field statistics and 928 effective properties of power-law 2D polycrystals (Lebensohn et al., 2004a, 929 2005) and 3D cubic, hexagonal (Lebensohn et al., 2004b) and orthorhombic 930 (Castelnau et al., 2008a) materials. The FFT-based formulation was also ap-931 plied to compute the development of local misorientations in polycrystalline

copper, with direct input from orientation images (Lebensohn et al., 2008). 933 As will be detailed is Section 4.2 it was further extended to transient behav-934 ior with an elasto-viscoplastic formulation (Idiart et al., 2006; Suguet et al., 935 2011; Lebensohn et al., 2012). The FFT-based full-field formulation for vis-936 coplastic polycrystals is conceived for a periodic unit cell, provides an exact 937 solution of the governing equations, and has better numerical performance 938 than a FE calculation for the same purpose and resolution. The viscoplas-939 tic FFT-based formulation consists in finding a strain-rate field, associated 940 with a kinematically-admissible velocity field, which minimizes the average 941 of local work-rate, under the compatibility and equilibrium constraints. The 942 method is based on the fact that the local mechanical response of a periodic 943 heterogeneous medium can be calculated as a convolution integral between 944 the Green function of a linear reference homogeneous medium and the actual heterogeneity field. Such type of integrals reduce to a simple product 946 in Fourier space, therefore the FFT algorithm can be used to transform the 947 heterogeneity field into Fourier space and, in turn, to get the mechanical fields by antitransforming that product back to real space. However, since 949 the actual heterogeneity field depends precisely on the a priori unknown me-950 chanical fields, an iterative scheme should be implemented to obtain, upon 951 convergence, a compatible strain-rate field and a stress field in equilibrium. 952 The periodic unit cell representing the polycrystal is discretized by means 953 of a regular grid  $\{x^d\}$ , which in turn determines a corresponding grid of the 954 same dimensions in Fourier space  $\{\xi^d\}$ . Velocities and tractions along the boundary of the unit cell are left undetermined under the sole condition of 956 periodicity. An average velocity gradient  $V_{i,j}$  is imposed to the unit cell, 957 which gives an average strain rate  $\dot{\bar{\varepsilon}}_{ij} = \frac{1}{2}(V_{i,j} + V_{j,i})$ . The local strain-rate field is a function of the local velocity field, i.e.  $\dot{\varepsilon}_{ij}(\mathbf{v}_k(\boldsymbol{x}))$ , and can be

split into its average and a fluctuation term:  $\dot{\varepsilon}_{ij}(\mathbf{v}_k(\boldsymbol{x})) = \dot{\varepsilon}_{ij} + \tilde{\varepsilon}_{ij}(\tilde{\mathbf{v}}_k(\boldsymbol{x})),$ where  $\mathbf{v}_i(\boldsymbol{x}) = \dot{\varepsilon}_{ij}x_j + \tilde{\mathbf{v}}_i(\boldsymbol{x})$ . By imposing periodic boundary conditions, the velocity fluctuation field  $\tilde{\mathbf{v}}_k(\boldsymbol{x})$  is assumed to be periodic across the boundary of the unit cell, while the traction field is antiperiodic, to meet equilibrium on the boundary between contiguous unit cells. The local constitutive equation that relates the deviatoric stress  $\boldsymbol{\sigma}'(\boldsymbol{x})$  and the strain rate  $\dot{\boldsymbol{\varepsilon}}(\boldsymbol{x})$  at point  $\boldsymbol{x}$ is obtained from Eqs (18) to (20).

If p(x) is the unknown pressure field introduced by the incompressibility constraint, the Cauchy stress field can be written as:

$$\sigma(x) = L^0 : \dot{\varepsilon}(x) + \varphi(x) - p(x)I$$
(26)

where the polarization field  $\varphi(x)$  is given by:

$$\varphi(\mathbf{x}) = \sigma'(\mathbf{x}) - \mathbf{L}^0 : \dot{\varepsilon}(\mathbf{x})$$
 (27)

where  $L^0$  is the stiffness (viscosity) of a linear reference medium. Eqs. (26) and (27) amount to transform the actual heterogeneity problem into an equivalent one, corresponding to a homogenous medium with eigen-strainrates. Note, however, that the above defined polarization field depends on the unknown  $\dot{\varepsilon}(\boldsymbol{x})$ . Combining Eq. (27) with the equilibrium and the incompressibity conditions gives:

$$L_{ijkl}^{0}v_{k,lj}(\boldsymbol{x}) + \boldsymbol{\varphi}_{ij,j}(\boldsymbol{x}) - p_{,i}(\boldsymbol{x}) = 0, v_{k,k}(\boldsymbol{x}) = 0$$
(28)

Assuming for a moment that the polarization field  $\varphi(x)$  is known, the system of partial differential equations (28), with periodic boundary conditions across the unit cell boundary, can be solved by means of the Green function method.

If  $G_{km}$  and  $H_m$  are the periodic Green functions associated with the velocity

and hydrostatic pressure fields, the solutions of system (28) are convolution integrals between those Green functions and the actual polarization term. The velocity gradient, after some manipulation is given by:

$$\tilde{v}_{i,j}(\boldsymbol{x}) = \int_{R^3} G_{ik,jl}(\boldsymbol{x} - \boldsymbol{x'}) \varphi_{kl}(\boldsymbol{x'}) d\boldsymbol{x'}.$$
 (29)

Convolution integrals in direct space are simply products in Fourier space.

where  $\hat{\Gamma}^{sym}_{ijkl} = \text{sym}(\hat{G}_{ik,jl})$ . The tensors  $\hat{G}_{ik}(\xi)$  and  $\hat{\Gamma}^{sym}_{ijkl}(\xi)$  are only func-

tions of  $L^0$  and can be readily obtained for every point belonging to  $\{\xi^d\}$ 

985 Hence:

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$$\hat{\dot{\tilde{\epsilon}}}_{ij}(\xi) = \hat{\Gamma}_{ijkl}^{sym}(\xi)\hat{\varphi}_{kl}(\xi), \tag{30}$$

(for details, see Lebensohn et al. (2008)). Now, taking into account the 988 definition 27 of  $\varphi(x)$ , Eq. 29 is an integral equation where the velocity 989 gradient appears in both sides, and, thus, it can be solved iteratively. Assigning initial guess values to the strain-rate field in the regular grid (e.g.  $\tilde{\dot{\epsilon}}$   $(x^d) = 0 \Rightarrow \dot{\epsilon}^{(0)}(x^d) = \dot{\bar{\epsilon}}$  ), and computing the corresponding stress field  $\sigma'^{(0)}(x^d)$  from the local constitutive relation (18) allows to obtain an initial guess for the polarization field in direct space  $\varphi^{(0)}(x^d)$  (27), which in turn 994 can be Fourier-transformed to obtain  $\hat{\boldsymbol{\varphi}}^{(0)}(\boldsymbol{\xi}^d)$ . 995 The rate of convergence of this fixed point technique is rather poor for 996 nonlinear constitutive relations such as power-law relations between the 997 stress and the strain-rate. Accelerated schemes based on augmented La-998 grangians have been proposed to improve this rate of convergence originally 999 by Michel et al. (2000, 2001) for composites, and later adapted by Lebensohn 1000 (2001) for polycrystals to which the interested reader in referred for details. 1001 Upon convergence, the stress at each material point can be used to calculate 1002 the shear rates associated with each slip system (Eq. 4), from which fields 1003 of relative activity of the basal, prismatic and pyramidal slip modes can be 1004

obtained, as well. While it is certainly possible to use the FFT-based for-1005 mulation for the prediction of microstructure evolution, in this section we 1006 have restricted our analysis to the local fields that are obtained for a fixed 1007 configuration. In this sense, the high strain-rate regions predicted by the 1008 model (see below) should be regarded as precursors of localization bands. 1009 Evidently, microstructural changes that are not considered under this ap-1010 proximation, like the eventual grain's and subgrain's morphologic evolution 1011 and rotation, as well as the possible occurrence of local strain hardening, 1012 may modify some of the trends observed in the initial micromechanical fields. 1013 In order to account for these microstructural changes, the FFT-based for-1014 mulation has been coupled with the front-tracking numerical platform Elle 1015 (Bons et al., 2001). Results of this coupled model are reported in Section 1016 5.2.2. 1017

1018 4.1.2. Application to columnar ice deforming in the secondary creep regime.

Lebensohn et al. (2009) and Montagnat et al. (2011) applied this FFT 1019 method to simulate strain rate and stress fields, and local lattice misorienta-1020 tions obtained at secondary creep in columnar ice polycrystals. Lebensohn 1021 et al. (2009) compared the simulated fields to a series of compression creep 1022 experiments performed by Mansuy et al. (2000, 2002) on laboratory-grown 1023 columnar ice samples characterized by multicrystals of controlled shape and 1024 orientations. The specimen used for this comparison (see Fig. 13) was a 1025 plate of 210×140 mm with a relatively thick (8 mm) section, consisting of 1026 a multicrystalline cluster, located in the center of the plate, with c-axes 1027 lying on the plane of the plate, and embedded in a matrix of fine-grained 1028 ice. This specimen was deformed under a compressive stress of 0.75 MPa 1029 exerted vertically in the plane at -10°C under plane strain conditions. Fig. 1030

13 shows, after 0.07 strain, three types of localization bands: basal shear 1031 bands, kink bands and sub-boundaries, that change orientation to follow 1032 crystallographic directions when they cross from one grain to another. 1033 In this configuration, kink band boundaries are seen mainly inside grains 1034 oriented close to 45° from the imposed compression direction. Kink bands, 1035 described as a sharp or discontinuous change in orientation of the active 1036 slip surface, had been reported in experimental studies conducted on 2-D 1037 ice polycrystals (Wilson et al., 1986; Wilson and Zhang, 1994; Montagnat 1038 et al., 2011). Sub-boundaries parallel to the c-axis were also observed. 1039

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The FFT-based calculation as described in the previous section was run to obtain the overall and local mechanical response of the above-described unit cell representing a columnar ice polycrystal, to the following imposed strain-rate tensor (see also Fig. 14):

$$\dot{\bar{\varepsilon}}_{ij} = \begin{bmatrix} 1 \times 10^{-8} & 0 & 0 \\ 0 & -1 \times 10^{-8} & 0 \\ 0 & 0 & 0 \end{bmatrix} s^{-1}$$
 (31)

The crystallographic texture of the 2-D ice polycrystal consisting of colum-1045 nar grains with c-axes perpendicular to the axial (vertical) direction x<sub>3</sub> was 1046 described in terms of a collection of Euler-angle triplets of the form  $(\varphi_1, 90^{\circ},$ 1047  $\varphi_2$ ) (Bunge convention). The application of the FFT method required the 1048 generation of a periodic unit cell or representative volume Element (RVE), 1049 by repetition along  $x_1$  and  $x_2$  of a square domain. This square domain was 1050 constructed in such a way that it contained the cross-sections of 200 colum-1051 nar grains, generated by Voronoi tessellation (see Fig. 14). This square do-1052 main is the cross-section of the unit cell, consisting of columnar grains with 1053

axes along  $x_3$  and sections in the  $x_1$ - $x_2$  plane. This unit cell was discretized 1054 using a  $1024 \times 1024 \times 1$  grid of regularly-spaced Fourier points, resulting in 1055 an average of around 5250 Fourier points per grain. Note that the periodic 1056 repetition of this unit cell along x<sub>3</sub> determines infinitely long grains along 1057 this direction. Three specific orientations with c-axis respectively at 0°, 45°, 1058 and 90° from the compression direction were forced to be among the set of 1059 200 (otherwise random) orientations assigned to the grains. For a plane-1060 strain state, such that  $x_1$  is the tensile direction and  $x_2$  is the compression 1061 direction, the grain with  $\varphi_1 = 45^{\circ}(45 \text{ deg grain in what follows})$  is theo-1062 retically favorably oriented to deform by soft basal slip, while in the 0 deg 1063 and 90 deg grains, the hard pyramidal systems are the only ones favorably 1064 oriented to accommodate deformation. It is worth noting that due to the 1065 above plane-strain condition and the in-plane orientation of the c-axes, the 1066 prismatic slip systems are not well-oriented, for any  $\varphi_1$  angle. 1067

The computed effective response of this kind of isotropic columnar ice 1068 polycrystal deformed in-plane is twice softer compared to an isotropic 3-D 1069 polycrystalline ice (Lebensohn et al., 2007). The computed overall relative 1070 activities of the different slip modes (i.e. 90.7%, 7.6% and 1.7% for basal, 1071 pyramidal and prismatic slip, respectively) show a preeminence of basal slip, 1072 a minor contribution of pyramidal slip and a very low activity of prismatic 1073 slip. Fig. 15 shows the computed equivalent strain-rate field for the entire 1074 unit cell, normalized with respect to the average equivalent strain rate ( $\dot{\bar{\varepsilon}}_{eq}$  = 1075  $1.15 \times 10^{-8} \text{s}^{-1}$ ). The main feature observed in this plot is a network of high 1076 strain-rate bands, precursors of localization bands (in what follows we will 1077 sometimes refer to them simply as "localization bands"). These bands are 1078 transmitted from grain to grain and are, in general, inclined with respect 1079 to the shortening and extension directions. They follow tortuous paths, 1080

sometimes with large deviations from  $\pm 45$  ° (i.e. the macroscopic directions 1081 of maximum shear stress). They follow crystallographic directions (basal 1082 poles or basal planes) inside each grain. The predicted bands parallel or 1083 perpendicular to the c-axis were reasonably assumed to be kink or shear 1084 bands, respectively (see Lebensohn et al. (2009) for details). Some segments 1085 of these bands also follow favorably-oriented grain boundaries and frequently 1086 go through triple or multiple points between grains, in good agreement with 1087 some of the observations of (Mansuy et al., 2002) (Fig. 13). Fig. 16 shows in 1088 more details the predicted fields of equivalent strain rate (normalized to  $\dot{\bar{\varepsilon}}_{eq}$ ), 1089 equivalent stress (in units of  $\tau^{bas}$ ) and relative basal activity, in the vicinities 1090 of the 45 deg grain. Two very intense (i.e. local strain rates higher than 10 1091 times the macroscopic strain rate) and parallel kink bands are seen inside 1092 the 45 deg grain, connected by several less intense shear bands (orthogonal 1093 to the pair of kink bands, lying on to the basal plane), in good agreement 1094 with Mansuy's experiments (see Fig. 13). The basal activity in the 45 deg 1095 grain is very high, although some regions of high non-basal activity can be 1096 observed between shear bands and immediately outside the kink bands. The 1097 latter is compatible with a low or even vanishing resolved shear stress on 1098 basal planes in those locations, which may be responsible for the formation 1099 of basal dislocation walls that are at the origin of a kink band (Mansuy 1100 et al., 2002). This correlation between kink band precursors and nearby 1101 localized high non-basal activity is systematic in these results. From the 1102 same detailed analysis performed around the 0 and 90 deg grains, a good 1103 match was found with experimental observations. 1104

In (Montagnat et al., 2011), the viscoplastic FFT-based approach was applied to the exact experimental microstructure of a compressive test performed on a 2D columnar sample. Samples (dimensions  $\approx 10 \times 10 \times 1.5$  cm<sup>3</sup>)

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were grown in the laboratory under a uniaxial temperature gradient to reach 1108 a columnar microstructure with all c-axes lying parallel to the sample sur-1109 face. In this work, the observed kink bands could be associated with mis-1110 orientations between adjacent regions of a grain interior of more than 5°, 1111 and their exact nature in term of dislocation arrangements were confirmed 1112 by EBSD measurements. Although the boundary conditions of the model-1113 ing were slightly different from the experimental one, the model was able 1114 to predict the exact location of the localization bands. The bands were as-1115 sociated with stress concentration that could reach five times the applied 1116 macroscopic stress, and to high levels of local non basal activity (see Fig. 1117 17). Nevertheless, the amplitude of the modeled lattice misorientations were 1118 always overestimated, and this was associated with the fact that very local 1119 grain boundary migration and new grain nucleation (dynamic recrystalliza-1120 tion mechanisms) observed experimentally were not considered in the model 1121 (see Section 5.2.2). 1122

## 1123 4.2. Elasto-viscoplastic FFT approach

The full-field FFT approach described above has been extended very re-1124 cently to the case of elasto-viscoplasticity, see (Idiart et al., 2006; Suguet 1125 et al., 2011; Lebensohn et al., 2012). As for purely viscoplastic behaviors, 1126 its application to highly anisotropic material like ice allows investigating the 1127 accuracy of elasto-viscoplastic mean-field models (see Section 3.6) since, as 1128 already mentioned, the FFT technique provides the "exact" (in a numerical 1129 sense) response of the specimen with the actual microstructure and local con-1130 stitutive relations. Application to ice allows studying transient creep effects 1131 with more detail. Comparison with experimental strain field measured with 1132 an intragranular spatial resolution has been provided in (Grennerat et al., 1133

2012) making use of the relative ease of producing samples with controlled 2-D microstructure, compared to other polycrystalline materials.

### 1136 4.2.1. The mechanical problem

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The method described in Suguet et al. (2011) considers the same mi-1137 crostructure description as in Section 4.1: a polycrystalline volume V com-1138 posed of several grains of different orientations, each grain obeying con-1139 stitutive relations defined in Section 2.3. The volume V is subjected to 1140 a macroscopic loading path, which can be a prescribed history of average 1141 strain, or a history of average stress or a combination of both. For simplicity, 1142 the method is presented here assuming a prescribed history of macroscopic 1143 strain  $\bar{\varepsilon}(t), t \in [0, T]$ . Other types of loadings can be handled by different 1144 methods described in (Michel et al., 1999) for instance. 1145

The local problem to be solved to determine the local stress and strain fields in the volume element V consists of the equilibrium equations, compatibility conditions, constitutive relations and periodicity boundary conditions:

$$\begin{cases} (\dot{\boldsymbol{\sigma}}, \dot{\boldsymbol{\tau}}_{0}, \dot{\boldsymbol{X}}) = \boldsymbol{F}(\dot{\boldsymbol{\varepsilon}}, \boldsymbol{\sigma}, \boldsymbol{\tau}_{0}, \boldsymbol{X}, \boldsymbol{x}, t), & \text{for } (\boldsymbol{x}, t) \in V \times [0, T], \\ \boldsymbol{\varepsilon}(\boldsymbol{x}, t) = \frac{1}{2} (\nabla \boldsymbol{u}(\boldsymbol{x}, t) + {}^{\mathsf{T}} \nabla \boldsymbol{u}(\boldsymbol{x}, t)), \\ \text{div } \boldsymbol{\sigma}(\boldsymbol{x}, t) = 0 & \text{for } (\boldsymbol{x}, t) \in V \times [0, T], \\ \boldsymbol{u}(\boldsymbol{x}, t) - \overline{\boldsymbol{\varepsilon}}(t).\boldsymbol{x} \text{ periodic on } \partial V, & \text{for } t \in [0, T] \end{cases}$$

$$(32)$$

The data of interest are the effective response  $\overline{\sigma}(t)$ ,  $t \in [0,T]$  of the polycrystal, the history of the average strain  $\overline{\varepsilon}(t)$ ,  $t \in [0,T]$ , but also the local
fields  $\sigma(x,t), \varepsilon(x,t)$  and other significant fields (internal variables, thermodynamic forces etc....).

The extension of the simplest version of the FFT-based method, also

called the *basic scheme*, to constitutive relations including crystalline elastoviscoplasticity relies on two ingredients:

1. A time-integration scheme for the constitutive differential equations. The time interval of interest [0,T] is split into time steps  $[t_n,t_{n+1}]$ . All quantities are assumed to be known at time  $t_n$ , and the quantities at time  $t_{n+1}$  are unknown. This time integration is performed at every point  $\mathbf{x}^d$  of the discretized polycrystal and the evolution problem is reduced to a problem for the stress and strain fields  $\boldsymbol{\sigma}$  and  $\boldsymbol{\varepsilon}$  at time  $t_{n+1}$  in the form

$$\sigma_{n+1}(\mathbf{x}^d) = \mathcal{F}_{n+1}(\mathbf{x}^d, \varepsilon_{n+1}(\mathbf{x}^d))$$
(33)

2. A FFT global scheme to solve the local problem for a nonlinear composite obeying Eq. (33).

The algorithm developed applies to a wide class of constitutive relations, see (Suquet et al., 2011). As before, it is limited to specimens submitted to periodic boundary conditions. Results presented below are performed with the FFT-based program Craft (freely available at http://craft.lma.cnrs-mrs.fr). For application to elastoviscoplasticity in ice, the local constitutive relation is the one provided above, see Eqs (2, 3, 6, 7, 8). It can also be formulated via the following differential equation:

$$\dot{\mathbf{Y}} = \mathbf{F}(\dot{\boldsymbol{\varepsilon}}, \mathbf{Y}, t), \tag{34}$$

1173 where

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$$oldsymbol{Y} = \left(egin{array}{ccc} oldsymbol{\sigma} \\ au_0^{(k)}, \ k=1,...M \end{array}
ight),$$

$$F(\dot{\boldsymbol{\varepsilon}}, \boldsymbol{Y}, t) = \begin{pmatrix} C : \left(\dot{\boldsymbol{\varepsilon}} - \sum_{k=1}^{M} \dot{\gamma}^{(k)}(\boldsymbol{Y})\boldsymbol{\mu}^{(k)}\right) \\ \left(\tau_{sta}^{(k)} - \tau_{0}^{(k)}\right) \sum_{\ell=1}^{M} h^{(k,\ell)} \left|\dot{\gamma}^{(\ell)}(\boldsymbol{Y})\right| \\ c^{(k)}\dot{\gamma}^{(k)}(\boldsymbol{Y}) - d^{(k)}X^{(k)} \left|\dot{\gamma}^{(k)}(\boldsymbol{Y})\right| - e^{(k)} \left|X^{(k)}\right|^{m} \operatorname{sign}\left(X^{(k)}\right) \end{pmatrix}$$
(35)

with  $m{C}$  the elastic stiffness ( $m{C}=m{S}^{-1}$ ). The set of parameters used are given in Table 1.

1176 4.2.2. Application to strain field prediction in a 2D-1/2 configuration.

The elasto-viscoplastic FFT approach was used to predict strain and 1177 stress field evolution during transient creep tests on ice polycrystals, in 1178 comparison with experimental measurements performed by Grennerat et al. 1179 (2012).1180 Samples were grown following (Montagnat et al., 2011) (see Section 4.1.2). 1181 This way, when compressed, (i) plastic deformation can be approximate as 2-1182 D, and (ii) strain fields measured at the specimen surface are representative 1183 for the sample volume owing to the minimisation of in-depth microstructure 1184 gradients. Average grain size (section perpendicular to the column direction) 1185

was about 5 mm and most of the c-axes were oriented parallel to the surface

 $(\pm 15^{\circ})$ . The microstructure and grain orientation were measured using an Automatic Fabric Analyzer (Russell-Head and Wilson, 2001) which provides orientation values with about 50  $\mu$ m resolution, and 1° accuracy. A Digital Image Correlation technique (Vacher et al., 1999) was applied to measure the strain heterogeneities on the surface perpendicular to the column direction. From displacement measurements performed during transient creep in ice, i.e. up to 1 to 2%, at -10°C, under 0.5 MPa, strain fields were evaluated with a resolution of about 0.2%, and at a spatial resolution of about 1 mm.

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The experimental microstructures were implemented in the code using the fabric analyzer data of  $2000\times2000$  pixels (but the model input does not need to be square). One pixel in the third dimension (column direction) is enough to reproduce the 2D-1/2 geometry thanks to the periodic boundary conditions.

Fig. 18 presents the strain field measured experimentally at the end of 1201 the transient creep, and the simulated fields of strain and stress. Although 1202 simulated boundary conditions did not precisely match the experimental 1203 ones, the heterogeneities of the strain field that develop during transient 1204 creep of polycrystalline ice were reproduced well by the model (Grennerat 1205 et al., 2012). In particular, the model was able to reproduce the characteris-1206 tic length of the heterogeneities being larger than the grain size, and scaling 1207 with the sample dimensions. Furthermore, both experimental and modeled 1208 results showed no correlation between the orientation of the c-axis and the 1209 strain intensity (see Fig. 19). This result casts doubt on the relevance of 1210 the distinction between "hard grains" and "soft grains" classically made for 1211 the analysis of ice mechanical behavior, and more generally for anisotropic 1212 materials. 1213

Fig. 20 represents the evolution of the simulated equivalent strain field from 0.25 to 0.60% of compression during transient creep. As observed experimentally, the strain heterogeneities develop early during the transient creep and are reinforced up to about 10 times the imposed strain.

## 5. Modeling of dynamic recrystallization mechanisms

Under laboratory conditions (described in Section 1), dynamic recrys-1219 tallization (DRX) dominates the changes of microstructures and fabrics in 1220 the tertiary creep regime, that is after about 1% macroscopic strain (Du-1221 val, 1981; Jacka and Maccagnan, 1984; Jacka and Li, 1994). During DRX, 1222 grain nucleation and grain boundary migration are two processes that con-1223 tribute to the reduction of the dislocation density, therefore of the stored 1224 deformation energy (Humphreys and Hatherly, 2004). In the laboratory, 1225 tertiary creep is a continuous sequence of deformation and recrystallization 1226 that gradually results in a steady state. This steady state is associated with 1227 an equilibrium grain size (Jacka and Li, 1994) and a girdle-type fabric with 1228 c-axes at about 30° from the compression axis (Jacka and Maccagnan, 1984), 1229 or with two maxima in simple shear (Bouchez and Duval, 1982). 1230 In polar ice sheets, DRX was identified from observation on ice thin sections 1231 along ice cores (Alley, 1992; Thorsteinsson et al., 1997; de la Chapelle et al., 1232 1998; Kipfstuhl et al., 2006). Three regions are usually defined: (i) normal 1233 grain growth driven by the reduction of grain-boundary energy in the upper 1234 hundreds meters of the core, (ii) rotation recrystallization during which new 1235 grains are formed by the progressive lattice rotation of the subgrains in the 1236 main part of the core and (iii) migration recrystallization similar to the one 1237 observed in the laboratory, in the bottom part where the temperature is 1238

above -10°C (see Montagnat et al. (2009) and Faria et al. (this issue) for a review).

Recrystallization and grain growth significantly influence the microstructure, the fabric and therefore the mechanical properties. To be able to integrate these mechanisms in the modeling of ice deformation is therefore crucial for an accurate prediction of its behavior.

## 5.1. Dynamic recrystallization within mean-field approaches

Several attempts were made to integrate dynamic recrystallization mechanisms into mean-field approaches as described in Section 3.

On the basis of the VPSC scheme (tangent version) described in Section 3 1248 for the description of the mechanical behavior, Wenk et al. (1997) developed 1249 a nucleation and grain-growth model to represent DRX in anisotropic ma-1250 terials such as ice. The model is based on the hypothesis that grains with 1251 a high stored energy (highly deformed) are likely to nucleate new grains 1252 and become dislocation-free. They may also be invaded by their neighbors 1253 which have a lower stored energy. Depending on the respective importance 1254 of nucleation and grain boundary migration processes, the recrystallization 1255 textures are expected to favor either highly deformed components or less 1256 deformed ones. 1257

One must first remember that, in the VPSC scheme, grains are represented by inclusions in an homogenous equivalent medium (HEM). Grain interactions are therefore represented "averaged" through the interaction between the inclusion and the HEM.

In this model, nucleation is represented, by a probability of nucleation Pper time increment  $\Delta t$  for each deformation step:

$$P \propto \Delta t \times exp(-A/E^2) \tag{36}$$

The constant A depends on the grain boundary energy and was taken as 1264 an adjustable parameter. E is a proxy of the stored energy,  $E \propto \sum_{s} (\Delta \tau_0^s)$ 1265 with  $\Delta \tau_0^s$  the variation of the critical resolved shear stress on the system s1266 during the deformation step. This calculation supposes a hardening law for 1267 each slip system to be defined. An isotropic hardening law was chosen in the 1268 form  $\dot{\tau}_0^s = H \sum_s \dot{\gamma}^s$ , with hardening matrix H being isotropic. A threshold 1269 was then defined for the minimum strain energy to nucleate, and the new 1270 grain completely replaced the old one (same size, same orientation), with a 1271 stored energy equal to zero. 1272

The grain boundary migration rate was taken proportional to the difference 1273 in stored energy between the grain and the average, i.e., the HEM. The de-1274 velopment of the microstructure is therefore a balance between nucleation 1275 and growth. Adjustable parameters were varied arbitrarily for comparison 1276 purpose. 1277

Applied to ice, this model resulted into weaker fabrics than the one ob-1278 tained by the classical VPSC tangent approach, mostly because grains near 1279 the compression axis disappeared (high stored energy) and only a few girdle 1280 grains, and a few grains exactly aligned with the compression axis from the 1281 beginning, remained. 1282

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Thorsteinsson (2002) included some DRX in its Nearest-Neighbor Inter-1284 action (NNI) model described in Section 3. Polygonization associated with 1285 rotation (or continuous) recrystallization is accounted for by comparing the resolved shear stress in the crystal  $(|\sum_s \tau_s \hat{\boldsymbol{b}}_s|)$  to the applied stress (with  $\tau_s$  the shear stress on system s, and  $\boldsymbol{b}_s$  a unit vector in the direction of the 1288 Burgers vector). If the ratio is smaller than a given value, and the dislocation density higher than a given value, then the crystal size is halved and 1290

both new grains are rotated by a fixed  $\Delta\theta$  of 5°. Grain growth occurs by 1291 normal grain growth according to (Gow, 1969; Alley et al., 1986) parabolic 1292 law  $(D^2 - D_0^2 = Kt)$ . The grain growth factor K follows an Arhenius-type 1293 dependence on the temperature. To take into account the grain growth as-1294 sociated with the difference in dislocation-stored energy between the grain i 1295 and the average, this growth factor was modified into  $(\tilde{K} = (E_{disl}^{av} - E_{disl}^{i})K')$ 1296 with K' a constant depending on temperature and impurities. 1297 Migration recrystallization is included in the model by considering the bal-1298 ance between grain-boundary energy, and stored energy associated with dis-1299 locations (the stored energy is calculated following (Wenk et al., 1997), as 1300 just described, and translated into dislocation density). A crystal recrystal-1301 lizes (i.e. is replaced by a crystal with initial dislocation density  $\rho_0$ ) when 1302 the dislocation energy is higher than the grain-boundary energy. This as-1303 sumption relies on the hypothesis that stored energy is released by normal 1304 grain growth (driven by GB energy), and that dynamic recrystallization only 1305 occurs if this relaxation in not efficient enough to decrease the dislocation 1306 density. The size of the new crystal is adjusted with the effective stress fol-1307 lowing (Guillopé and Poirier, 1979; Jacka and Li, 1994) and its orientation 1308 is chosen at random in the range of the "softest" orientations in the applied 1309 stress state. 1310 Modeling results were obtained for comparison to a case similar to the GRIP 1311 ice core, with vertical compression, and rotation recrystallization domi-1312 nating. The introduction of polygonization allows for the preferential re-1313 moval of "hard" grains, which leads to a weaker fabric compared to the 1314 "no-recrystallization" case. In particular, when associated with the NNI 1315 formulation, the model is able to reproduce fabrics quite similar to those 1316 measured along the GRIP ice core at several depths. "Girdle-type" fabric 1317

similar to the experimental fabrics, results from the introduction of migration recrystallization. However, parametrization remains weak, in particular the estimation of the dislocation density, and of the recrystallized grain orientations.

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The last example presented here is the cellular automaton model for 1323 fabric development by Ktitarev et al. (2002) and Faria et al. (2002). The 1324 application was mostly to reproduce the fabric measured along deep ice 1325 cores, with the assumption of deformation under uniaxial compression. The 1326 cellular automaton (CA) frame is especially suitable for simulation of sys-1327 tems represented by a certain number of cells, which are associated with 1328 generalized state variables and arranged in regular environment. The con-1329 sidered material is a thin horizontal layer of ice located along the ice core, 1330 thin enough so that it is considered homogeneous in the vertical dimension. 1331 To discretize the problem according to the CA method, the authors took a 1332 one-dimensional lattice of equal cells representing the grains, described by 1333 their size, and their orientation. The basic dynamical quantity of the al-1334 gorithm is the dislocation density. This density increases with deformation 1335 and depends on the orientation of the grain. Recrystallization mechanisms 1336 proceed when a critical value is reached. Normal grain growth is accounted 1337 for following Gow (1969) and is apparently the only growth mechanism asso-1338 ciated with polygonization mechanisms. The increase in dislocation density 1339 is associated with the resolved shear stress on the basal system and the re-1340 crystallization model developed in Montagnat and Duval (2000) is used to 1341 estimate the evolution of the density in relation with grain size and polygo-1342 nization mechanisms. Rotation of grains is ruled by a kinematic equation 1343 based on the inelastic spin, assuming a compressive stress proportional to the 1344

depth along the core, and a linear dependence between the shearing rate of 1345 sliding on the basal system and the resolved shear stress. The time evolution 1346 was related to the depth along the core using the Dansgaard et al. (1993) 1347 relation. Following Duval and Castelnau (1995), migration recrystallization 1348 was only applied bellow 2800 m depth. During migration recrystallization, 1349 new grains were allowed to grow much faster by consuming up to ten cells 1350 at every time step, until it is impinged by another growing grain, or until it 1351 reaches the critical size of the steady state. 1352 The model was able to provide a good qualitative evolution of the grain 1353 size, by separating the influence of normal grain growth, polygonization and 1354 migration recrystallization similarly to what was suggested from the mea-1355 surements along the GRIP ice core (Thorsteinsson et al., 1997; de la Chapelle 1356 et al., 1998). Concerning the fabric evolution, the model was able to pre-1357 dict the evolution toward a single maximum, but the kinetics is too strongly 1358 influenced by the polygonization, and further by migration recrystallization. 1359

#### 1361 5.2. Dynamic recrystallization within full-field approaches

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This section presents a coupling between a platform for structural change 1362 in materials (Elle) with the full-field FFT approach presented in Section 4.1, 1363 to predict the microstructure evolution of ice polycrystals during dynamic 1364 recrystallization. A critical step in the development of generic models linking 1365 plastic deformation and recrystallization is the incorporation of the interac-1366 tion between intra- and intergranular heterogeneities of the micromechanical 1367 fields (i.e. strain rate and stress) and the recrystallization processes. Be-1368 cause local rotations of the crystal lattice are controlled by local gradients of 1369 plastic deformation, heterogeneous distributions of lattice orientations are 1370

observed at the grain and subgrain scale, see Section 4. This has a strong 1371 influence on recrystallization as this is a process driven by the local gradi-1372 ents of energy (e.g. grain boundary or stored strain energy). Traditional 1373 mean-field models used to predict microstructure evolution during recrys-1374 tallization are based on a simplified description of the medium and cannot 1375 fully describe intragranular heterogeneities (Section 5.1). Therefore, explicit 1376 full-field approaches are required for a better understanding of dynamic re-1377 crystallization and prediction of microstructure evolution at large strain. 1378

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## 1380 5.2.1. The Elle modeling plateform

Elle is a platform for the numerical simulation of processes in rocks and grain aggregates, with particular focus on (micro-) structural changes (Jessell et al., 2001; Jessell and Bons, 2002; Bons et al., 2008; Piazolo et al., 2010). The simulations act on an actual 2D image of the microstructure (Fig. 21.a). Elle is currently restricted to 2D cases although the underlying principles for 2D are equally valid in 3D (Becker et al., 2008), and therefore the approach could be converted for 3D simulations.

The central philosophy of Elle is to enable the coupling of processes 1388 that act on the material, recognizing that the effect of one process may 1389 significantly alter that of a concurrent process. Dynamic recrystallization, 1390 for example, can greatly change crystallographic preferred orientations in 1391 mineral aggregates deforming by dislocation creep (Jessell, 1988a,b). Cou-1392 pling of processes is achieved in Elle using the principle of operator splitting, 1393 whereby individual processes successively act on the model in isolation, for 1394 a small time step. This approach greatly simplifies coding, as the coupling 1395 between processes needs not be programmed itself, but emerges from their 1396

1397 alternating effect on the model.

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Each process in Elle is an individual program or module. A shell-script takes the starting model and then passes it in a loop to the individual processes, which each in turn modify the model slightly. Each loop represents one time step. The user can freely determine the mix of processes that operate by choosing which ones to include in the loop. The relative activity of individual processes is determined by the parameters passed on to each process.

The model is essentially defined by two types of nodes: boundary nodes 1405 (bnodes) and unconnected nodes (unodes) (Fig. 21.b). Bnodes define the 1406 boundaries of a contiguous set of polygons (termed flynns). These flynns 1407 typically represent single grains, but can also represent regions within a 1408 material, for example rock layers (Llorens et al., 2012). The boundaries of 1409 the flynns are formed by straight segments that connect neighboring bnodes. 1410 One brode can be connected by either two or three other brodes. The use 1411 of bnodes and flynns makes the model suitable for a range of Finite Element and front-tracking models. 1413

Unodes form a second layer of the model. These are nodes that do not necessarily have fixed neighborhood relationships and typically represent points within the material. Some processes are not amenable to be modeled with polygons, but are best simulated with a regular grid of unodes. The FFT code is an example. Nodes and flynns can have a range of attributes assigned to them, such as c-axis orientation, boundary properties, etc.

Elle uses fully wrapping boundaries. A flynn that touches one side of the model continues on the other side (Fig. 21.a). The model is thus effectively a unit cell that is repeated infinitely in all directions. Although Elle typically uses a square model, deformation may change the unit cell into a 1424 parallelogram shape.

Elle now includes a large and ever growing number of process modules for a variety of processes that mostly relate to microstructural developments in mineral aggregates. Each process can essentially act on the model in only two ways: changing the position of a node (e.g. a bnode in case of grain boundary migration) or changing the value(s) of attributes of flynns or nodes (e.g. concentration at a unode in a diffusion simulation). Some of the most relevant current processes are:

• Normal grain growth driven by the reduction of surface energy, and hence curvature of grain boundaries. This process was used by Roessiger et al. (2011) to address the issue of the competition between grain growth and grain size reduction in the upper levels of polar ice caps (Mathiesen et al., 2004). Surface energy can be anisotropic, i.e. depending on the lattice orientation of the grains on either side of the boundary (Bons et al., 2001). Two-phase grain growth has been applied to grain growth in rocks with a small proportion of melt (Becker et al., 2008) and to ice with air bubbles (Fig. 21)(Roessiger et al. this volume).

• The Finite Element module Basil is used for incompressible power-law viscous deformation (Barr and Houseman, 1996). Using viscosities that are assigned to flynns, it calculates the stress and velocity fields resulting from applied boundary conditions. It has been used to study the behaviour of rigid inclusions in a deforming matrix (Bons et al., 1997), the behaviour of deforming two-phase materials as a function of viscosity contrast and composition (Jessell et al., 2009) and for folding

of layers (Llorens et al., 2012). The wrapping boundaries of the Elle model, in combination with continuous remeshing allows for arbitrarily high strains (Jessell et al., 2009). In combination with grain growth and dynamic viscosity, Jessell et al. (2005) studied strain localisation behaviour. Durand et al. (2004) investigated the influence of uniaxial deformation on grain size evolution in polar ice cores and its influence on ice dating methods.

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- Dynamic recrystallisation includes grain boundary migration driven by strain energy (dislocation density) and the formation of new grain boundaries by progressive subgrain rotation or polygonisation (Urai et al., 1986). In the next section (5.2.2) we will describe how these processes, employing a front-tracking model for grain-boundary migration, are linked with the FFT approach (Griera et al., 2011, 2012; Piazolo et al., 2012) to model the stress and strain-rate fields and the driving forces for recrystallization.
- A final Elle module of potential relevance to ice is that developed by Schmatz (2010) for the interaction between migrating grain boundaries and small particles (e.g. dust or clathrates). The particles are represented by unodes, which, when swept by a grain boundary, can latch onto that boundary. Particles can slow down grain boundary movement, but can also be dragged along and eventually be released by a grain boundary.

Summarizing, Elle provides a large number of routines to simulate grain-1472 scale processes in minerals and rocks, and hence in glacial or polar ice. The open and versatile code allows for more process modules to be added or existing ones to be tailored for application to ice. A significant advantage of 1475

the code is that it enables the investigation of the complex microstructural and mechanical effects of multiple, concurrent and coupled processes.

### 1478 5.2.2. Coupling Elle platform to FFT approach

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Most of the numerical approaches used to simulate deformation and mi-1479 crostructural evolution of rocks and metals are based on combining defor-1480 mation approaches based on the Finite Element Method with Monte Carlo, 1481 cellular automaton, phase field, network or level-set methods to simulate 1482 recrystallization (Jessell, 1988a,b; Raabe and Becker, 2000; Piazolo et al., 1483 2002, 2010, 2012; Solas et al., 2004; Battaile et al., 2007; Logé et al., 2008). 1484 An alternative to these methods is the numerical scheme used in this study 1485 based on the coupling between the crystal plasticity FFT-based code (Leben-1486 sohn, 2001) (section 4.1) and the Elle modeling platform just described 1487 (Bons et al., 2008). Both codes have been previously explained and here 1488 we only concentrate on some particularities of the coupling between them. 1489 The FFT-based formulation is integrated within the Elle plateform using 1490 a direct one-to-one mapping between data structures. The polycrystalline 1491 aggregate is discretized into a periodic, regular array of spaced and uncon-1492 nected nodes (Fourier Points in the FFT and "unconnected nodes", unodes, 1493 in Elle; Fig. 21). 1494

Numerical simulation is achieved by iterative application of small time steps of each process. After numerical convergence of the FFT model, data is transferred to Elle assuming that the micromechanical fields are constant in the incremental time step. The position and material information of unodes are directly updated because they are equivalent to the Fourier points, while position of boundary nodes (bnodes) are calculated using the velocity field. Based on the evolution of the predicted local lattice rotation field, the dislo-

cation density can be estimated using strain gradient plasticity theory (e.g. 1502 Gao et al. (1999); Brinckmann et al. (2006)) or using the dislocation density 1503 tensor or Nyes tensor (Nye, 1953; Arsenlis and Parks, 1999; Pantleon, 2008). 1504 With this approach, only geometrical necessary dislocations required to en-1505 sure strain compatibility are estimated. To simplify the problem, we use 1506 a scalar approach where all dislocations are assumed to be related to the 1507 basal plane. The lattice-orientation and dislocation-density fields provide 1508 the input parameters to predict recrystallization in the aggregate. 1500

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Recrystallization is simulated by means of three main processes: nucle-1510 ation, subgrain rotation and grain boundary migration. Using the kinematic and thermodynamic instability criteria of classical recrystallization theory 1512 (Humphreys and Hatherly, 2004; Raabe and Becker, 2000), nucleation is 1513 simulated by the creation of a small new, dislocation-free flynn when the local misorientation or dislocation density exceeds a defined threshold. The 1515 lattice orientation of the new grain is set to that of the critical unode. When 1516 a cluster of unodes within a grain share the same orientation that is different from the rest of the unodes in that grain, a new grain boundary is created, 1518 while preserving the lattice orientations of the unodes. A technical limita-1519 tion is that nucleation of grains and subgrains is only allowed along grain boundaries. Nucleation within grains are therefore not possible.

Grain boundary migration is described by a linear relationship between 1522 velocity (v) and driving force per unit area ( $\Delta f$ ), by  $v = M\Delta f$  where M 1523 corresponds to the grain boundary mobility, which has an Arhenius-type dependency on temperature. Grain boundary curvature and stored strain energies are used as driving forces for grain boundary motion. For this

situation, the driving force can be defined as

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$$\Delta f = \Delta E - 2\gamma/r \tag{37}$$

is the boundary energy and r is the local radius of curvature of the grain 1529 boundary. Stored strain energy is the energy per unit volume associated 1530 with lattice distortions and depends on the dislocation density  $(\rho)$  and dis-1531 location type. Grain boundary motion is simulated using the free-energy 1532 minimization front-tracking scheme of (Becker et al., 2008). When an un-1533 ode is swept by a moving grain boundary, it is assumed that dislocations 1534 are removed and the new lattice orientation is that of the nearest unode 1535 belonging to the growing grain. 1536 Following the Elle philosophy, each process runs individually, following a 1537 pre-established sequence. After all Elle processes have run, the unodes layer 1538 is used to define the new input microstructure to be deformed viscoplasti-1539 cally by the FFT code. A drawback is that the unodes are not following a 1540 regular mesh, a requirement needed by the FFT approach. For this reason, 1541 as proposed by Lahellec et al. (2003), and later adapted in the context of 1542 Elle by Griera et al. (2011, 2012), a particle-in-cell method is used to remap 1543 all material and morphological information to a new regular computational 1544 mesh. In order to avoid unrealistic crystallographic orientations, these are 1545 not interpolated during remapping. The crystallographic orientation of a 1546 new Fourier Point that belongs to a specific grain is that of the nearest un-1547 ode that belongs to the same grain. This allows to run numerical simulation 1548 up to large strains. 1549

where  $\Delta E$  is the difference of stored strain energy across the boundary,  $\gamma$ 

### 1550 5.2.3. Application to creep experiments and natural ice

An example of numerical simulation using the FFT/Elle approach is 1551 shown in Fig. 22. The simulation is based on a creep experiment of poly-1552 crystalline columnar ice. Samples and experimental conditions are those 1553 of (Montagnat et al., 2011) described in Section 4.1.2. The specimen was 1554 deformed at -10°C under uniaxial conditions with a constant load of 0.5 1555 MPa up to an axial strain of 4%. A thin section of the initial and the 1556 final microstructure was analyzed using the Automatic Ice Texture Ana-1557 lyzer method (Russell-Head and Wilson, 2001) to obtain the local c-axis 1558 orientations. After a 4% of shortening, the onset of local recrystallization 1559 is evident in the experiment (Fig. 22a), in the form of irregular and ser-1560 rated grain boundaries and small new grains that are preferentially located 1561 at triple junctions and along grain boundaries. Localized variations in the 1562 orientation of the basal plane form sharp and straight subgrain boundaries 1563 that indicate intracrystalline deformation. The experimental c-axis map was 1564 used as input for the FFT/Elle simulation. The experimental starting mi-1565 crostructure was discretized into a grid of 256×256 Fourier points. As only 1566 the c-axis orientation is known, the other axes are given a random orienta-1567 tion. Crystal plasticity is described with an incompressible rate-dependent 1568 equation for basal, prismatic and pyramidal slip (see Section 3). Critical 1569 resolved shear stress for basal slip was set 20 times lower than for non-basal 1570 systems. The physical properties used for recrystallization are as follows: 1571 mobility  $M=1\times 10^{-10}~\mathrm{m^2Kg^{-1}s^{-1}}$  (e.g. Nasello et al., 2005), isotropic 1572 boundary energy  $\gamma = 0.065~\mathrm{Jm^{-2}}$  (Ketcham and Hobbs, 1969), shear mod-1573 ulus  $G = 3 \times 10^9$  Pa and critical dislocation density  $\rho = 1 \times 10^{12} \text{ m}^{-2}$  (de la 1574 Chapelle et al., 1998). Pure shear boundary conditions were imposed with 1575

vertical constant strain rate of  $-1 \times 10^{-8}$  s<sup>-1</sup> up to a 4% of strain in 1% increments.

The computed orientation map and grain boundary misorientation are shown in Fig. 22b. Several features of the experiment are seen in the numerical simulation, such as the development of sharp misorientations or kink bands, bulging and serrated grain boundaries, and new grains at triple junction and grain boundaries. There is a good correlation between loca-tion of kink bands in the experiment and the simulation. However, the width of kink bands in the simulation is dependent on the numerical reso-lution. A relationship between grain boundary motion/nucleation and high dislocation-density regions is observed (Fig. 23). Variations in dislocation densities across grain boundaries lead to migration of these boundaries in the direction of the dislocation density gradient. However, some discrepan-cies are also seen, such as, for example, grain boundary motion (e.g. at the bottom-left part) that was not observed in the experiment. One explanation may be that low and high angle grain boundaries were not differentiated in the simulation and, therefore, both types had similar mobility. 

A second example aims to show the strong effect recrystallization can have on the final microstructure. A  $10\times10~\mathrm{cm^2}$  microstructure with 1600 grains with random c-axis orientations (Fig. 24a) was deformed to 40% shortening in plane-strain pure shear. The values of mechanical (slip systems, CRSS, etc) and recrystallization (mobility, surface energy, etc) properties are similar to those of the model described before, but adjusted to a natural strain rate of  $10^{-12}~\mathrm{s^{-1}}$  at about -30°C. Fig. 24b shows the c-axis and relative misorientation maps for an extreme case with no recrystallization (FFT only). Dominant red and purple colors indicate that the c-axis of crys-

tallites are preferentially oriented at low angles to the shortening direction. 1603 Elongated grains are oriented parallel to the stretching direction. Remark-1604 able differences are observed when recrystallization is activated (Fig. 24c). 1605 Grain boundaries are smooth and grains larger and more equidimensional. 1606 Despite the significant difference in microstructure, both simulations show a 1607 single maximum c-axis distribution at low angle to the shortening direction. 1608 The strong resemblance of the simulated microstructure with that of natural 1609 ice (Thorsteinsson et al., 1997; de la Chapelle et al., 1998; Weikusat et al., 1610 2009) shows the strong potential of modeling of ice deformation based on 1611 an actual map of the microstructure. 1612

#### 6. Toward large scale ice flow modeling 1613

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A number of models have been developed in glaciology to simulate the 1614 flow of anisotropic ice and the strain-induced development of fabric within 1615 polar ice-sheets. Accounting for ice anisotropy in an ice-flow model implies 1616 to (i) build a macroscopic anisotropic flow law whose response will depend on the local fabric and (ii) have a proper description of the ice fabric at 1618 each node of the mesh domain and be able to model the fabric evolution as 1619 a function of the flow conditions. We hereafter present the main issues to 1620 address these two points.

Due to the scale of these large ice-masses, the implementation of a polycrystalline law must stay simple enough and numerically tractable. At present, full-field or even homogenization models presented previously are computationally too demanding and cannot realistically be used to estimate the mechanical response in an ice-sheet flow model. Here we present two approaches to build a simple and efficient macroscopic law for polycrystalline ice. The first one is based on the concept of a scalar enhancement factor function so that the collinearity between the strain-rate and the deviatoric stress tensors is conserved (Placidi and Hutter, 2006), see Section 6.1. The second polycrystalline law is fully orthotropic and depends on six relative viscosities, function of the fabric (Gillet-Chaulet et al., 2005, 2006), see Section 6.2. Both models are phenomenological and must be calibrated using experimental or numerical results, as described below.

With regards to other materials, the advantage of the hexagonal symme-1635 try of ice is that the crystal rheology can be assumed transversally isotropic 1636 (only true for a linear rheology). Under this assumption, only one unit vector 1637 suffices to describe the lattice orientation, thus simplifying the mathemat-1638 ical description of fabrics. With regards to other materials, the advantage 1639 of the hexagonal symmetry of ice is that only one unit vector suffices to de-1640 scribe the lattice orientation, thus simplifying the mathematical description 1641 of fabrics. The discrete description of the fabric, i.e. a couple of angles for 1642 each crystal, would require too large a number of variables to be stored at 1643 each node of the domain mesh. Typical mesh size are hundreds of thousand 1644 nodes in 3D (Seddik et al., 2011) up to few millions for the most recent appli-1645 cations (Gillet-Chaulet et al., 2012). The use of a parameterized orientation 1646 distribution function (ODF) would decrease the number of parameters, but 1647 evolution equations for these parameters to describe the fabric evolution 1648 cannot be obtained in a general case (Gagliardini et al., 2009). The ori-1649 entation tensors, which describe the fabric at the macroscopic scale in a 1650 condensed way are more suitable. Five parameters are needed to describe 1651 an orthotropic fabric (the two eigenvalues of the second-order orientation 1652 tensor and the three Euler angles to specify the position of the material 1653 symmetry basis), and an evolution equation for the second-order orienta-1654

tion tensor can be easily derived from the macroscopic stress and strain-rate fields.

#### 6.1. Continuous Diversity and the CAFFE model

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The CAFFE model (Continuum-mechanical Anisotropic Flow model based on an anisotropic Flow Enhancement factor) results from a suitable combination of two basic concepts: a power law description of ice rheology resembling the well-known Glen's flow law (Glen, 1955); and a multiscale approach to model the evolution of the polycrystalline microstructure of ice based on the general theory of continuous diversity (Faria, 2001; Faria and Hutter, 2002; Faria et al., 2003).

The ideas leading to the CAFFE model have been elaborated in a series 1665 of works by Luca Placidi and his collaborators (Placidi, 2004, 2005; Placidi 1666 and Hutter, 2005, 2006; Placidi et al., 2004). These ideas culminated in the 1667 definitive CAFFE formulation, presented by Placidi et al. (2010), in which 1668 the so-called enhancement factor of Glen's flow law becomes a function of 1669 the material anisotropy (fabric), and the evolution of the latter is governed 1670 by an orientation-dependent mass balance equation derived from the theory 1671 of continuous diversity applied to glacier and ice-sheet dynamics (Faria, 1672 2006a,b; Faria et al., 2006). 1673

The greatest strength of the CAFFE model is its successful compromise between accuracy and flexibility, which allows one to upgrade existing computer models of isotropic ice-sheet dynamics based on Glen's flow law into efficient anisotropic models, without profound changes in the original code. In fact, due to its relative simplicity, the CAFFE model has already been implemented in several numerical ice-flow simulations. For instance, it has been used by Seddik et al. (2008) and Bargmann et al. (2011) to simulate

the ice flow at the site of the EPICA-DML drill site at Kohnen Station,
Dronning Maud Land, East Antarctica, while Seddik et al. (2011) used it
to simulate the ice flow in the vicinity of the Dome Fuji drill site in central
East Antarctica.

In the following, we review the CAFFE formulation presented by Placidi et al. (2010). The fundamental idea is to regard polycrystalline ice as a "mixture" of lattice orientations, following the philosophy of the theory of Mixtures with Continuous Diversity (MCD) proposed by Faria (2001, 2006a). Succinctly, a mixture with continuous diversity is a multicomponent medium made up of an infinite number of mutually interacting species, whose distinctive properties vary smoothly from one to another.

In the case of polycrystalline ice, species are distinguished by their c-axis 1692 orientations. Each point of the continuous body is interpreted as a representative volume element, which encompasses a large number of crystallites with 1694 their own c-axis orientations. Each of such orientations is mathematically 1695 identified with a point on the surface of the unit sphere S<sup>2</sup> and represented by a unit vector  $n \in \mathbb{S}^2$ . As a consequence, for each species one can intro-1697 duce a mass density field  $\rho^*(\boldsymbol{x},t,\boldsymbol{n})$ , given at a certain position  $\boldsymbol{x}$  within the 1698 polycrystal, and at time t, sometimes called orientational mass density, such 1699 that, when integrated over the whole unit sphere, the usual mass density 1700 field of the polycristal (i.e. of the "mixture") results: 1701

$$\varrho(\boldsymbol{x},t) = \int_{\mathbb{S}^2} \varrho^*(\boldsymbol{x},t,\boldsymbol{n}) \,\mathrm{d}^2 n , \qquad (38)$$

where  $d^2n$  (=  $\sin\theta d\theta d\phi$  in spherical coordinates) is the infinitesimal solid angle on the unit sphere  $S^2$ . The product  $\varrho^*(\boldsymbol{x},t,\boldsymbol{n}) d^2n$  is the mass fraction of crystalline material in the volume element with c-axis directed towards  $\boldsymbol{n}$  within the solid angle  $d^2n$ . Therefore, assuming that the material is incompressible, the mass (or volume) fraction  $\varrho^*/\varrho$  can be interpreted as the usual orientation distribution function (ODF) in the context of materials science (Bunge, 1993; Zhang and Jenkins, 1993; Raabe and Roters, 2004). It should be remarked that in the glaciological literature the term "ODF" sometimes refers to the relative number, instead of the mass (or volume) fraction, of grains with a certain orientation.

The time evolution of  $\varrho^*$  is governed by the balance equation of species (orientational) mass

$$\frac{\partial \varrho^*}{\partial t} + \operatorname{div}(\varrho^* \boldsymbol{v}) + \operatorname{div}_{\boldsymbol{n}}(\varrho^* \boldsymbol{u}^*) = \varrho^* \Gamma^*$$
(39)

1714 with

$$\operatorname{div}_{n}(\boldsymbol{\Phi}^{*}) = \operatorname{tr}\left[\operatorname{grad}_{n}(\boldsymbol{\Phi}^{*})\right],$$

$$\operatorname{grad}_{n}(\boldsymbol{\Phi}^{*}) = \frac{\partial \boldsymbol{\Phi}^{*}}{\partial n} - \left(\frac{\partial \boldsymbol{\Phi}^{*}}{\partial n} \cdot n\right) n$$
(40)

for any scalar-, vector- or tensor-valued field  $\Phi^*(\boldsymbol{x},t,\boldsymbol{n})$ . In (39),  $\boldsymbol{u}^*(\boldsymbol{x},t,\boldsymbol{n})$  denotes a sort of "velocity" on the unit sphere (with  $\boldsymbol{u}^*\cdot\boldsymbol{n}=0$ ), called orientational transition rate. Further,  $\Gamma^*(\boldsymbol{x},t,\boldsymbol{n})$  is the specific recrystallization rate, which describes the rate of change of mass (per unit mass) of one species into another one with different orientation. Integration of (39) over the unit sphere  $S^2$  gives rise to the usual mass balance equation for the polycrystal (i.e. the "mixture")

$$\frac{\partial \varrho}{\partial t} + \operatorname{div}(\varrho \boldsymbol{v}) = 0 \quad \text{with}$$

$$\int_{\mathbb{S}^2} \varrho^* \Gamma^* \, \mathrm{d}^2 n = \int_{\mathbb{S}^2} \operatorname{div}_{\boldsymbol{n}} \left(\varrho^* \boldsymbol{u}^*\right) \, \mathrm{d}^2 n = 0,$$
(41)

Notice that the first integral in (41) is a consequence of mass conservation, while the second integral follows from Gauss' theorem.

As shown by Faria (2001, 2006a) and Faria and Hutter (2002), the transition rate  $u^*$  is governed by its own balance equation, involving couple stresses and body couples. In the development of the CAFFE model, however, an abridged approach has been adopted by postulating a constitutive equation for the transition rate

$$\boldsymbol{u}^* = \boldsymbol{W}\boldsymbol{n} - \iota \left[ \dot{\boldsymbol{\varepsilon}}\boldsymbol{n} - (\boldsymbol{n} \cdot \dot{\boldsymbol{\varepsilon}}\boldsymbol{n}) \, \boldsymbol{n} \right] - \frac{\lambda}{\rho^*} \operatorname{grad}_{\boldsymbol{n}} \left( \varrho^* H^* \right)$$
(42)

1729 where

$$\boldsymbol{W} = \frac{1}{2} \left( \operatorname{grad} \boldsymbol{v} - (\operatorname{grad} \boldsymbol{v})^T \right), \dot{\boldsymbol{\varepsilon}} = \frac{1}{2} \left( \operatorname{grad} \boldsymbol{v} + (\operatorname{grad} \boldsymbol{v})^T \right)$$
(43)

are the tensors of rotation and strain rate, respectively. The first term on 1730 the right hand side of (42) represents a rigid-body rotation, while the second 1731 term describes the process of strain-induced lattice rotation (Dafalias, 2001), 1732 with  $\iota > 0$  denoting the so-called "shape factor" of the theory of rotational 1733 diffusion (Faria, 2001). According to Placidi et al. (2010), fabric evolution 1734 simulations of the GRIP and EPICA-DML ice cores suggest that best results 1735 are obtained for  $0.6 > \iota > 0.4$ . Finally, the third term on the right hand 1736 side of (42) models rotation recrystallization as a diffusive process, with 1737  $\lambda > 0$  being the orientational diffusivity and  $H^*(\boldsymbol{x},t,\boldsymbol{n})$  an orientational 1738 ("chemical") potential, also called "hardness function" by Gödert (2003). In 1739 principle  $H^*$  should be a constitutive function, but, based on microstructural 1740 analyses of the NorthGRIP ice core (Durand et al., 2008), Placidi et al. 1741 (2010) suggest that one may simply set  $H^* = 1$ . 1742 In the original application of the MCD theory to the flow of glaciers and 1743 ice sheets (Faria, 2006b), the specific recrystallization rate  $\Gamma^*$  is regarded as 1744 a dissipative variable. However, for simplicity, in the CAFFE model Placidi 1745 (2004, 2005) has proposed the following relation between  $\Gamma^*$  and the strain 1746

1747 rate

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$$\Gamma^* = G(D^* - D) , \text{ with}$$

$$D^* = 5 \frac{(\dot{\boldsymbol{\varepsilon}} \boldsymbol{n})^2 - (\boldsymbol{n} \cdot \dot{\boldsymbol{\varepsilon}} \boldsymbol{n})^2}{\operatorname{tr}(\dot{\boldsymbol{\varepsilon}}^2)} \text{ and } D = \frac{1}{\varrho} \int_{\mathbb{S}^2} \varrho^* D^* d^2 n ,$$
(44)

where G > 0 is a material parameter, while  $5/2 \ge D^* \ge 0$  and  $5/2 \ge D \ge 0$  are called the species and polycrystal "deformability", respectively.

As remarked by Placidi et al. (2010), owing to the difficulties in determining the values of the material parameters  $\lambda$  and G from experiments, they are usually determined by fitting numerical simulations of ice core fabrics and grain stereology. This concludes the description of the fabric evolution.

As for the flow law, in contrast to the full stress–strain rate relation with tensorial fluidity (viscosity) predicted by the theory of continuous diversity (Faria, 2006b), the CAFFE model adopts a much simplified generalization of Glen's flow law:

$$\dot{\boldsymbol{\varepsilon}} = E(D) A(T) \sigma_{eg}^{n-1} \boldsymbol{\sigma}' , \qquad (45)$$

where  $\boldsymbol{\sigma'}$  is the deviatoric part of the Cauchy stress tensor  $\boldsymbol{\sigma},~\sigma_{eq}$  is the 1758 effective stress invariant, n is the power law exponent (usually set equal 1759 3), T is the temperature, and A(T) is a temperature-dependent rate factor. 1760 Clearly, (45) implies that all anisotropy effects are contained in the scalar-1761 valued, deformability-dependent flow enhancement factor E(D), such that 1762 stress and strain rate are collinear and (45) reduces to the classical form of 1763 Glen's flow law when  $E(D) \equiv \text{const.}$ 1764 A detailed functional form for the enhancement factor E(D) has been 1765 proposed by Seddik et al. (2008) and Placidi et al. (2010), which is con-1766

tinuously differentiable at D=1 and is compatible with the experimental

results of Azuma (1995) and Miyamoto (1999)

$$E(D) = \begin{cases} (1 - E_{\min}) D^{\zeta} + E_{\min} & 1 \ge D \ge 0, \\ \frac{4D^{2} (E_{\max} - 1) + 25 - 4E_{\max}}{21} & 5/2 \ge D > 1, \end{cases}$$
(46)

1769 with

$$\zeta = \frac{8}{21} \left( \frac{E_{\text{max}} - 1}{1 - E_{\text{min}}} \right) , \qquad E_{\text{max}} \approx 10 , \quad E_{\text{min}} \approx 0.1 .$$
 (47)

By introducing the orientation tensors (essentially equivalent to the dipole and quadrupole moments of  $\varrho^*/\varrho$ )

$$a^{(2)} = \frac{1}{\varrho} \int_{\mathbb{S}^2} \varrho^* \boldsymbol{n} \otimes \boldsymbol{n} d^2 n, \quad \boldsymbol{a}^{(4)} = \frac{1}{\varrho} \int_{\mathbb{S}^2} \varrho^* \boldsymbol{n} \otimes \boldsymbol{n} \otimes \boldsymbol{n} \otimes \boldsymbol{n} d^2 n$$
 (48)

to reformulate the CAFFE flow law (45) in an explicitly anisotropic form

$$\dot{\boldsymbol{\varepsilon}} = \hat{E}(\boldsymbol{\sigma}') A(T) \sigma_{eq}^{n-1} \boldsymbol{\tau} , \qquad (49)$$

In plain words, (49) tells us that the CAFFE model can be applied to all anisotropies (fabrics) that can satisfactorily be represented by a multipole expansion up to fourth order. Fortunately, most anisotropies observed in glaciers and ice sheets.

## 1777 6.2. GOLF law and Elmer/Ice

In this section, we present the anisotropic ice flow model developed at LGGE. This model has been used for various applications (Gillet-Chaulet et al., 2005, 2006; Durand et al., 2007; Martín et al., 2009; Ma et al., 2010). In this approach, the fabric is described using the second and fourth-order orientation tensors (48). In this continuum description of the fabric, the polycrystal represents the local behavior of a representative elementary ice volume. By assuming that the fourth-order orientation tensor  $\boldsymbol{a}^{(4)}$  is given

as a tensorial function of  $a^{(2)}$  (Gillet-Chaulet et al., 2005), the fabric can 1785 be described in a very condensed way using  $a^{(2)}$  solely. By definition, 1786  ${\rm tr}\, {m a}^{(2)}=1,$  so that only the first two eigenvalues  $a_1^{(2)}$  and  $a_2^{(2)}$  and three 1787 Euler angles are needed to completely define the fabric. As a consequence, 1788 modeled fabrics are orthotropic, i.e. the c-axes distribution presents three 1789 orthogonal symmetry planes. Although orthotropy is a simple form of the 1790 most general anisotropy, it is thought to be a good compromise between 1791 physical adequateness and simplicity. The second-order orientation tensor 1792 allows to describe all the observed fabric patterns: for random c-axes dis-1793 tribution the diagonal entries of  $a^{(2)}$  are  $a_{11}^{(2)} = a_{22}^{(2)} = a_{33}^{(2)} = 1/3$ , for a 1794 single maximum fabric with its maximum in the third direction,  $a_{33}^{(2)}>1/3$ 1795 and  $a_{11}^{(2)} \approx a_{22}^{(2)} < 1/3$ , and for a girdle type fabric in the plane  $(x_1, x_2)$ , 1796  $a_{33}^{(2)}<1/3$  and  $a_{11}^{(2)}\approx a_{22}^{(2)}>1/3$ . When the material symmetry axes are 1797 those of the general reference frame, as for the three particular previous 1798 fabrics, the non-diagonal entries of  $a^{(2)}$  are zero. 1799

The behavior of the polycrystal is described by the general orthotropic 1800 linear flow law (GOLF, Gillet-Chaulet et al., 2005). In its initial form, ice 1801 was assumed to behave as a linearly viscous orthotropic material. In more 1802 recent works (Martín et al., 2009; Ma et al., 2010), the GOLF law has been 1803 extended to a nonlinear form by adding an invariant in the anisotropic linear 1804 law. The simple choice is either to add the second invariant of the strain 1805 rate (Martín et al., 2009) or the second invariant of the deviatoric stress 1806 (Pettit et al., 2007). No theoretical or experimental results are available 1807 today to discard one of these two solutions, and other solutions based on 1808 anisotropic invariants of the deviatoric stress and/or the strain rate are 1809 also possible. In (Ma et al., 2010) approach, the nonlinearity of the law is 1810 introduced through the second invariant of the deviatoric stress. With this 1811

definition, the anisotropy factors of the polycrystalline law for a given stress are identical in the linear and nonlinear cases. In other words, for a given fabric and a given state of stress, the corresponding strain rate relative to the isotropic response is the same for the linear and nonlinear cases. Using the strain-rate invariant in the same way as Martín et al. (2009) did, leads to different anisotropy factors (as defined here) in the linear and nonlinear cases. Therefore, the proposed expression of the nonlinear GOLF law is as follows:

$$\sum_{r=1}^{3} \left[ \eta_r tr(\boldsymbol{M}_r \cdot \dot{\boldsymbol{\varepsilon}}) \boldsymbol{M}'_r + \eta_{r+3} (\dot{\boldsymbol{\varepsilon}} \cdot \boldsymbol{M}_r + \boldsymbol{M}_r \cdot \dot{\boldsymbol{\varepsilon}})' \right] = 2A \sigma_{eq}^{n-1} \boldsymbol{\sigma}', \qquad (50)$$

where A is the temperature-dependent Glen's law parameter for isotropic ice. 1820 The six dimensionless anisotropy viscosities  $\eta_r(\boldsymbol{a}^{(2)})$  and  $\eta_{r+3}(\boldsymbol{a}^{(2)})$  (r=1,1821 (2,3) are functions of eigenvalues of the second-order orientation tensor  $a^{(2)}$ , 1822 which represent a measure of the anisotropy strength. The three structure 1823 tensors  $M_r$  are given by the dyadic products of the three eigenvectors of  $a^{(2)}$ , 1824 which then represent the material symmetry axes. In the method proposed by Gillet-Chaulet et al. (2005), the six dimensionless viscosities  $\eta_r(\boldsymbol{a}^{(2)})$  are 1826 tabulated as a function of the fabric strength  $(i.e., the a_i^{(2)})$  using a micro-1827 macro model. When ice is isotropic,  $\eta_r = 0$  and  $\eta_{r+3} = 1$  (r = 1, 2, 3), and 1828 Eq. (50) reduces to the isotropic Glen's flow law. 1829 Following Gillet-Chaulet et al. (2005), the six dimensionless viscosities 1830  $\eta_r(\boldsymbol{a}^{(2)})$  are tabulated using the visco-plastic self-consistent model (VPSC, 1831 Castelnau et al., 1996a, 1998), see Section 3. The two crystal parameters 1832 in the VPSC model used to tabulate the GOLF law were chosen so that 1833 the experimentally observed polycrystal anisotropy is reproduced. Gillet-1834 Chaulet et al. (2005) use the shear-strain rates ratio for a polycrystal with 1835 a single maximum fabric and an isotropic polycrystal both experiencing 1836

the same shear stress. This anisotropy factor in shear is hereafter noted 1837  $k_s$  and, according to the experimental results of Pimienta et al. (1987), its 1838 value is approximately  $k_s = 10$ . In other words, the VPSC parameters 1839 are chosen so that the response under simple shear of a polycrystal with a 1840 single maximum fabric is  $k_s$  times easier to deform than the corresponding 1841 isotropic polycrystal. The experimental results of Pimienta et al. (1987) also 1842 indicate that an isotropic polycrystal is much easier to deform than a single 1843 maximum fabric polycrystal experiencing the same uniaxial compressional 1844 These experiments allow to define a second anisotropy factor for 1845 uniaxial compressional stress, which is noted  $k_c$ . A value  $k_c = 0.4$  is in 1846 accordance with the experimental results of Pimienta et al. (1987). As 1847 discussed before, the anisotropy factors  $k_s$  and  $k_c$  are independent of Glen's 1848 flow law exponent n with the adopted nonlinear formulation. 1849

Assuming that recrystallization processes do not occur and that the ice fabric is induced solely by deformation, the evolution of the second-order orientation tensor  $a^{(2)}$  can be written as

$$\frac{Da^{(2)}}{Dt} = W \cdot a^{(2)} - a^{(2)} \cdot W - (C \cdot a^{(2)} + a^{(2)} \cdot C) + 2a^{(4)} : C, \quad (51)$$

where  $m{W}$  is the spin tensor defined as the antisymmetric part of the velocity gradient. The tensor  $m{C}$  is defined as

$$C = (1 - \alpha)\dot{\varepsilon} + \alpha k_s A \sigma_{eq}^{n-1} \sigma'.$$
 (52)

The interaction parameter  $\alpha$  controls the relative weighting of the strain rate  $\dot{\varepsilon}$  and the deviatoric stress  $\sigma'$  in the fabric evolution Eq. (51). When  $\alpha = 0$ , the fabric evolution is solely controlled by the state of strain rate, whereas in the case where  $\alpha = 1$  the fabric evolves under the influence of the deviatoric stress solely. In between, as for the VPSC, both the strain rate and deviatoric stress contribute to the fabric evolution. In what follows, the interaction parameter is  $\alpha = 0.06$  in accordance with the crystal anisotropy and the VPSC model used to derive the polycrystal behaviour (Gillet-Chaulet et al., 2005). In Eq. (51), the fourth-order orientation tensor is evaluated assuming a closure approximation giving  $\boldsymbol{a}^{(4)}$  as a tensorial function of  $\boldsymbol{a}^{(2)}$  (Gillet-Chaulet et al., 2005).

The anisotropic polycrystalline law described above and the associated 1866 fabric evolution equations have been implemented in the Finite Element 1867 code Elmer/Ice, the glaciological part of the open source Finite Element 1868 software Elmer developed by CSC (http://www.elmerfem.org/). Ice flow 1869 (velocity and isotropic pressure) are obtained solving the anisotropic Stokes 1870 equations and coupled with the fabric evolution equation (51) and the upper 1871 free surface equation in the case of transient simulations. In Gillet-Chaulet 1872 et al. (2006), the model was applied to synthetic geometries in order to 1873 show the influence of coupling the Stokes and fabric evolution equations on 1874 the flow of ice over a bumpy bedrock. In Durand et al. (2007), the model 1875 was used to explain the fabric evolution in the Dome C ice core, in the 1876 framework of the EPICA project. The authors showed that to explain the 1877 fabric evolution at Dome C, shear stress must be invoked. The model was 1878 also applied to evaluate the value of the ad-hoc enhancement factor that 1879 should be incorporated in large-scale isotropic ice-sheet flow model in Ma 1880 et al. (2010). In Martín et al. (2009), the anisotropic ice flow model was 1881 applied to explain observed shapes of isochrones below ridges or domes. 1882

## 7. Synthesis and perspectives

Applications of ice mechanical behavior modeling extend from below the single-crystal scale to the ice sheet scale. Upwards, this scale range far exceeds that of engineering material sciences but is similar to the geological one. Within this scale range, many physical processes come into play, some of which are not yet very well described. Furthermore, there exist strong interactions between these processes that create bridges between the different levels of complexity. Modeling of ice has strongly benefited from advances in materials science. In return, as shown by the results presented in this paper, the contribution of the ice community to the theoretical understanding and modeling of the mechanical behavior of anisotropic materials is significant. With the large viscoplastic anisotropy of the ice crystal, ice is now considered a model material. The advances presented here may equally well be applied to, for example, mantle flow, where the anisotropy due to fabric (CPO) development in olivine is thought to play a significant role (Tommasi et al., 2009; Long and Becker, 2010).

The presented modeling methods are basically of two types; some that aim to precisely reproduce the physical mechanisms as observed experimentally, and some with a more phenomenological approach. Going through scales, it clearly appears that individual dislocation interactions cannot be taken into account at the scale of the polycrystal imbedded in a glacier environment. Nevertheless, modeling at the scale of dislocation interactions provides a better estimate of the interactions between slip systems at the single crystal scale, which, in turn, is essential to reproduce an accurate mechanical response of the polycrystal with mean-field and full-field approaches. Furthermore, full-field approaches are necessary to validate the approximations

made using mean-field models, as they provide the "exact" (in a numerical sense) response of the specimen with a real microstructure, integrating the inter- and intra-granular interactions. Finally, large-scale flow models are now getting to a sufficient level of complexity to be able to take into account and represent the anisotropy associated with the fabrics induced by the flow conditions. To do so, they integrate mean-field approaches that correctly reproduce the viscoplastic anisotropy and a non-linear mechanical behavior.

A summary of the main domains of application, advantages, and limitations of the main modeling tools presented in the paper is given on tables 2, 3 and 4

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Much progress has recently been made in the modeling of dynamic re-1919 crystallization processes and their interactions with flow anisotropy. Never-1920 theless, due to the complexity of the physical processes involved, to jump 1921 the gap between scales is a strong challenge. The field dislocation mechanics 1922 approach appears very promising to associate the internal stress field and 1923 dislocation arrangements to the nucleation and grain boundary migration 1924 mechanisms. However, field dislocation mechanics cannot yet be applied to 1925 scales larger than the polycrystal. Full-field models, including the FFT-Elle 1926 coupling, have the same scale limitation, but may play an important role 1927 in parameterizing small-scale processes (dislocation glide, grain boundary 1928 migration, etc.) for mean-field models. They are also important tools to 1929 test models of mechanical and microstructural evolution. 1930

Compared to other minerals, ice shows remarkably strong transient behavior (Duval et al., 1983; Castelnau et al., 2008b). Continuum flow models, such as Glen law (Glen, 1955) have so far not been able to incorporate the resulting mechanical complexity of polycrystalline ice deformation. Only recently have mechanical models reached a level of sophistication to address

transient behavior. This development is promising and probably highly relevant in cases where ice flow changes at rates for where both elastic and
viscoplastic behavior may interact. In particular, this concerns the very
topical subject of ice shelves, ice streams or extra-terrestrial ice submitted
to tide forcing. Which model will be able to correctly take into account
these transient, and event cyclic behavior, and at which scale?

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A next step will likely be the multi-scale coupling of models of increasing complexity. We can expect dislocation dynamics and field dislocation mechanics to provide the local criteria for slip system interactions, nucleation, grain boundary migration as local input to full-field approaches that will be further used in interaction with mean-field approaches to calibrate dynamic recrystallization variables influencing the mechanical response and fabric development.

An interesting example of such model interweaving is given by the large-1949 scale flow modeling presented in this paper. Nevertheless, a strong effort 1950 is still required concerning the flow law of ice and its dependency on fab-1951 rics (CPO) and strain. Recent velocity measurements in Greenland (Gillet-1952 Chaulet et al., 2011) questioned the relevance of a stress exponent equal 1953 to three as classically considered for large scale flow modeling (for instance 1954 Paterson (1994); Hooke (2005); Greve and Blatter (2009), ...). Owing to 1955 the variety of processes that accommodate strain along an ice core path, 1956 one could also expect several regimes to occur with depth, as suggested by 1957 some authors (see for instance Lipenkov et al. (1989); Faria et al. (2009); 1958 Pettit et al. (2011)). Such modeling - observation comparisons mainly raise 1959 the complexity of the physical processes involved that can probably not be 1960 summarized in a single universal law.

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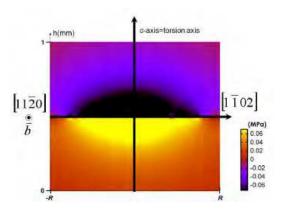


Figure 1: Map of the resolved shear stress in the prismatic system for a torsion boundary made of basal screw dislocations. The cylinder diameter is 1 mm, and the maximum applied stress is 0.1 MPa. From (Chevy et al., 2012)

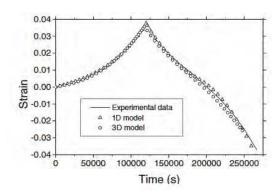


Figure 2: Creep curves in forward and reverse torsion from experiments on single crystals, obtained by 1D and 3D FDM models. From (Taupin et al., 2007)

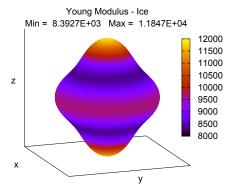


Figure 3: Young's modulus in [MPa] of an ice single crystal with its c-axis aligned with z, at  $-16^{\circ}{\rm C}$ .

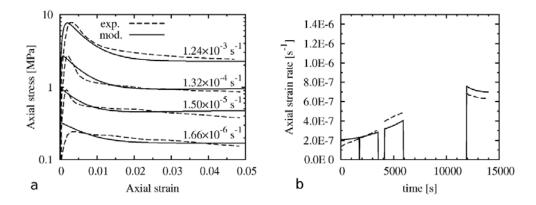


Figure 4: Behavior of an ice single crystal deformed by basal slip. a) Predictions of the model, based on Eq. (6) and with parameters given in Table 1, in comparison with the experimental data of Weertman (1973). Axial strain rates are indicated. b) Prediction of the model in comparison with the results of the recovery tests of Taupin et al. (2008). Temperature is  $-10^{\circ}$ C. From (Suquet et al., 2011)

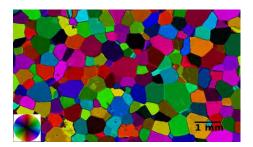


Figure 5: Typical 2D microstructure of an ice polycrystal grown in the laboratory. The color wheel gives the color-code for the c-axis orientation.

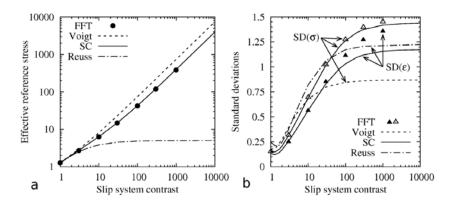


Figure 6: Full-field vs. mean-field behavior for ice polycrystals with random fabric, for a linear viscous behavior (n=1) and various viscoplastic anisotropy (or slip system contrats) at the grains level. a) Effective flow stress  $\tilde{\sigma}_0$ . b) Standard deviation of equivalent stress and strain rate, normalized by  $\bar{\sigma}_{eq}$  and  $\dot{\bar{\epsilon}}_{eq}$ , respectively, characterizing field heterogeneities at the polycrystal scale. Results from the linear SC scheme are compared to reference numerical solutions provided by the FFT approach. Reuss and Voigt bounds are also indicated. Note that, for these bounds, standard deviations of stress and strain rate, respectively, do vanish.

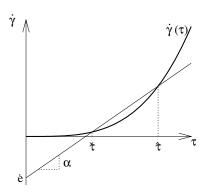


Figure 7: Schematic representation of the linearization between the shear rate  $(\dot{\gamma})$  and the stress  $(\tau)$ , to illustrate Eq. (21).

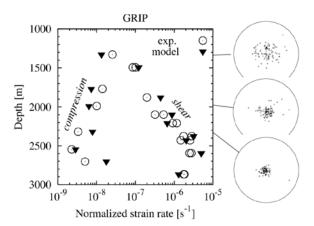


Figure 8: Stationary creep behavior at  $-10^{\circ}$ C calculated by the affine SC model, and compared to experimental data obtained on anisotropic specimens from the GRIP ice core. The c-axis pole figures on the right show an increasing concentration of c-axes towards the  $in\ situ$  vertical direction from the surface of the ice sheet down to  $\sim 2600$ m depth. Experimental data from Castelnau et al. (1998) are expressed for a stress of 1MPa using a stress sensitivity n=3. Points on the left hand side reflect the (hard) behavior under vertical compression, whereas data on the right correspond to (soft) horizontal shear.

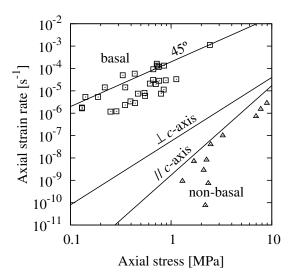


Figure 9: Stationary creep behavior of single crystals at  $-10^{\circ}$ C input in the AFF SC model to get results of Fig. 8 (lines), compared to the data set compiled by Duval et al. (1983) (symbols). Results are indicated for uniaxial compression at 45° from the c-axis (activation of basal slip), as well as for compression perpendicular (activation of prismatic slip) and parallel (activation of pyramidal systems) to the c-axis. From (Castelnau et al., 2008b)

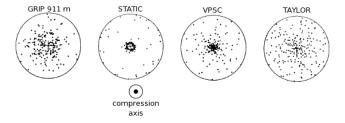


Figure 10: Comparison between fabrics measured along the GRIP ice core (911 m depth), simulated by the static (Reuss), VPSC-tangent, and Taylor (Voigt) approaches.

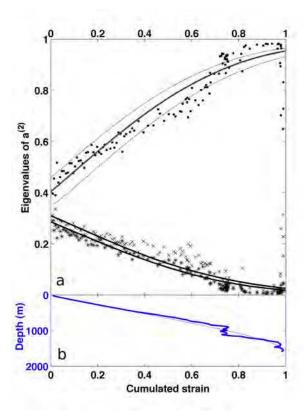


Figure 11: a) Evolution of the eigenvalues of the orientation tensor  $\mathbf{a}^{(2)} = \mathbf{c} \otimes \mathbf{c}$  of the fabric along the Talos Dome ice core, as a function of the cumulated compressive strain. Lines = VP-SO model results, dashed line represents the range of fabric evolution modeled with variation of the initial orientation tensor eigenvalue from isotropic (bottom line), as measured at 18 m (central line), more concentrated than measured (top line). Dots, crosses and plus = measurements performed with the Automatic Ice Texture Analyzer (Russell-Head and Wilson, 2001). b) Cumulated in-situ compressive strain as a function of depth as model by the TALDICE-1 chronology (full line) (Buiron et al., 2011).

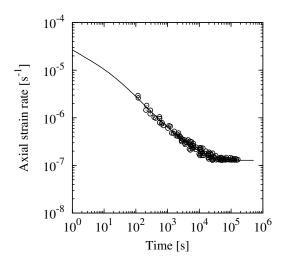


Figure 12: Transient creep response of isotropic ice under an uniaxial compressive stress of 1 MPa predicted with the affine elasto-viscoplastic extension of the self-consistent scheme (line). Model results are compared to the data of Ashby and Duval (1985), expressed for the same loading conditions (points). Strain hardening of prismatic and pyramidal slip systems is taken into account. From (Castelnau et al., 2008b)

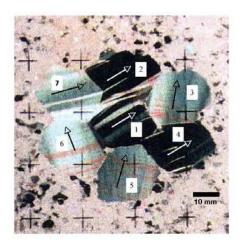


Figure 13: Photograph of a compression creep specimen (after (Mansuy et al., 2000)) between crossed polarizers, after a deformation of  $6.6 \times 10^{-2}$  at  $-10^{\circ}$ C. The corresponding strain rate was  $6.0 \times 10^{-8}$  s<sup>-1</sup>. The compression direction in vertical in the plane of the photograph. The mean size of each hexagonal grain was 20 mm. Black and white arrows indicate the initial c-axis orientations. Kink bands appear as abrupt changes in color parallel to the c-axis, shear bands are perpendicular to the c-axis direction.

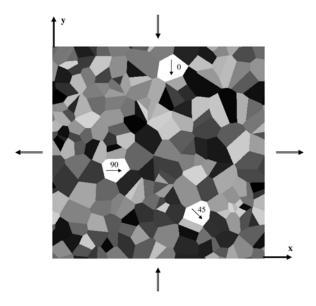


Figure 14: Unit cell containing the cross-sections of 200 columnar grains generated by Voronoi tesellation. The three hand-picked orientations:  $(0^{\circ}, 90^{\circ}, 0^{\circ})$ ,  $(45^{\circ}, 90^{\circ}, 0^{\circ})$  and  $(90^{\circ}, 90^{\circ}, 0^{\circ})$ , and the extension and shorting directions are also indicated.

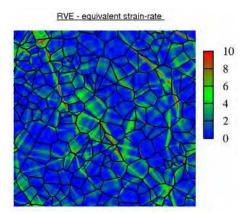


Figure 15: Predicted equivalent strain-rate field over the entire unit cell of Fig. 14, normalized with respect to the average equivalent strain rate ( $\dot{\bar{\varepsilon}}_{eq} = 1.15 \times 10^{-8} \mathrm{s}^{-1}$ )..

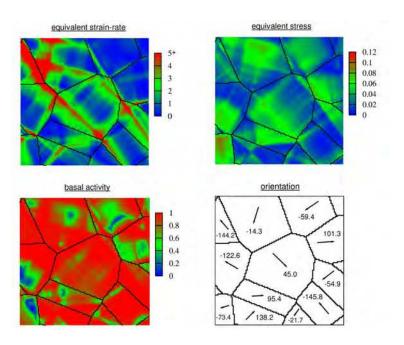


Figure 16: Predicted fields of equivalent strain rate (normalized to  $\dot{\bar{\varepsilon}}_{eq}$ ), equivalent stress (in units of  $\tau^{bas}$ ), relative basal activity, and map of neighbor orientations, for the 45 deg grain and its surroundings.

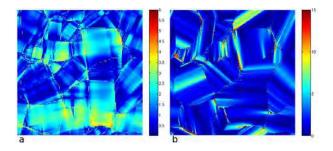


Figure 17: a) Predicted equivalent stress field (in units of  $\tau^{bas}$ ), and b) the misorientation, compared with initial orientation, obtained after 1% strain in a laboratory made microstructure. From (Montagnat et al., 2011)

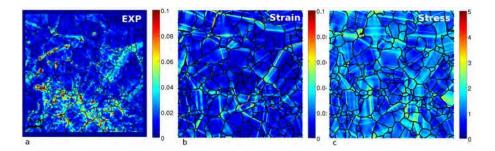


Figure 18: a) Strain field measured experimentally, b) simulated, and c) stress field simulated, after 0.85 % of axial compression. Experimental resolution is about  $75\times75$  pixels, the modeling one is  $1024\times1024$  pixels. From (Grennerat et al., 2012)

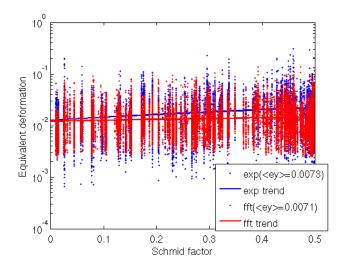


Figure 19: Equivalent strain as a function of the Schmid factor (as a proxy of the orientation). Experimental results are in blue, modeling results in red. Each point is one pixel of the microstructure. The macroscopic strain was 0.7%. From (Grennerat et al., 2012)

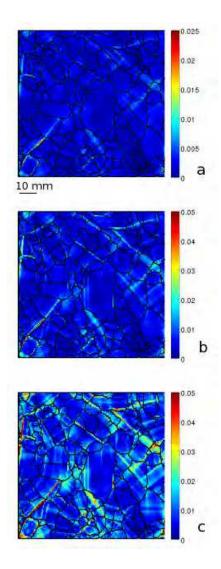


Figure 20: Evolution of the FFT-simulated equivalent strain field during the transient creep of a "2D-1/2" sample of ice, after (a) 0.15%, (b) 0.35% and (c) 0.60% compressive strain (see Grennerat et al. (2012)).

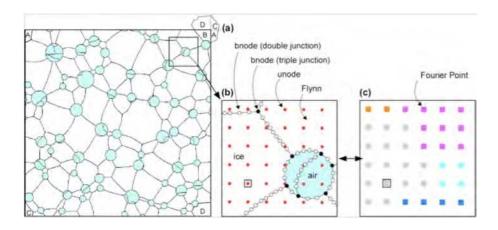


Figure 21: (a) Example of an Elle model: ice (white) with air bubbles (pale blue) (Roessiger et al., this volume). The Elle model has fully wrapping boundaries and grains A to D are in fact one single grain. (b) Close-up showing that grains are defined by flynns (polygons), themselves defined by straight segments that connect boundary nodes (bnodes). A second layer of unconnected nodes (unodes) can be added to keep track of material points. (c) For the FFT module, the microstructure is discretized into a periodic, regular mesh of Fourier Points defined by a characteristic lattice orientation. A direct mapping between unodes layer and Fourier points is established between both codes.

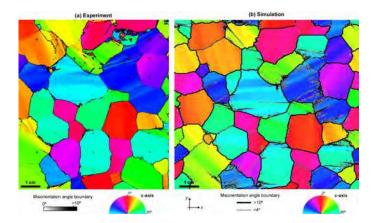


Figure 22: Comparison of (a) physical experiment and (b) numerical simulation after a vertical shortening of 4%. A qualitative equivalence between experiment and simulations is observable, such as correspondence of kink-bands or discontinuous subgrain boundaries at sharp grain boundaries asperities. Colors indicate the orientation of the c-axis respect to the sample reference. Misorientation angle between nodes are indicated in grey ( $>4^{\circ}$ ) and black ( $>10^{\circ}$ ). Triangular patches seen in the experiment are due to erroneous misfit during Automatic Ice Texture Analyzer acquisition.

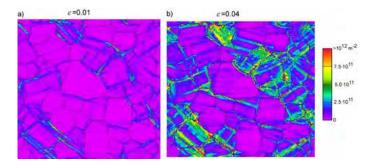


Figure 23: Dislocation density maps after (a) 1% and (b) 4% of shortening. Grain microstructure is indicated by dark lines. Serrated and bulging grain boundaries develop due to grain boundary migration into regions of high dislocation density. New recrystallized grains develop preferentially at triple points and along grain boundaries. Low dislocation densities are typically observed at bulge areas and new grains.

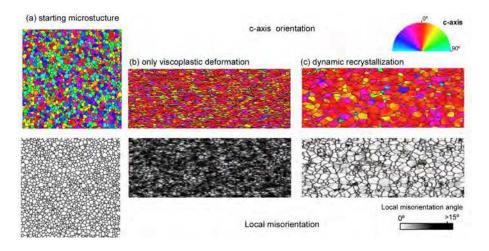


Figure 24: Numerical simulation of polar ice microstructure using the FFT/Elle scheme. (a) Starting microstructure. (b) 40% vertical shortening with only viscoplastic deformation. (c) 40% shortening with viscoplastic deformation coupled with recrystallization. Top row shows c-axis orientations in color and local misorientation in grey. Bottom row shows local misorientation only. C-axis orientation distributions are shown in lower-hemisphere stereoplots.

	$ au_{ini}$	$ au_{sta}$	$\dot{\gamma}_0$	n	c	d	e
Basal	0.1	0.022	$10^{-6}$	2	9	60	0.0003
Prismatic	0.13	1.5	$10^{-6}$	2.85	9	60	0.0003
Pyramidal	3.875	3.875	$10^{-6}$	4	9	60	0.0003

## Hardening matrix:

	Basal	Prismatic	Pyramidal
Basal	70	125	0
Prismatic	125	110	0
Pyramidal	0	0	0

Table 1: Material parameters used in the full-field simulations for single crystals of ice at  $-10^{\circ}C$ . Units are MPa and s<sup>-1</sup>.

Model	Scale	Type	Applications	Advantages	Limitations
DD	Single crystal (meso- scopic)	Finite Element (FE)	Dislocation dynamics and interactions	Solve physics of plasticity at the dislocation scale - Provide slip system activities and interactions	Computing time-intensive.  Small samples. Simplified configurations
FDM	Single to poly-crystal	Full-field (FE)	Plasticity modeling taking into account heterogeneous internal stress field associated with dislocations	Length scaling, makes the link between the complexity of the dislocation field dynamics and the macroscopic behavior - Provide intra-crystalline fields	Computing time-intensive - Limited number of grains
VPSC	Polycrystal	Mean-field	Provides effective visco-plastic behavior of polycrystals, based on a given single-crystal behavior	Texture-induced polycrystal anisotropy, texture evolution for secondary creep, can reach large strains	Very limited information on intra-crystalline fields, elasticity and recrystallization neglected
EVPSC	Polycrystal	Mean-field	Provides effective elasto-visco-plastic behavior at polycrystal level, based on known single-crystal behavior	Captures texture-induced anisotropy during transient creep regime	Very limited information on intra-crystalline fields, limited to small strains

Table 2: Summary of techniques application domains, interest and limitations. DD: Dislocation Dynamics, FDM: Field Dislocation Mechanics, VPSC: Homogenized Polycrystal Visco-Plasticity, EVPSC: Homogenized Polycrystal Elasto-Visco-Plasticity

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Model	Scale	Type	Applications	Advantages	Limitations
VPSC + DRX	Polycrystal	Mean-field	DRX mechanisms with phenomenological laws, fabric evolution	Fast to run for high strain - easily adaptable to various DRX laws	Too simplified description of DRX mechanisms. No account for intra-crystalline fields
VPFFT	Polycrystal	Full-field	Provides effective visco-plastic behavior and local intra- crystalline fields, 2D and 3D	Microstructural effects on local fields distribution in the secondary creep regime	Microstructure evolution at large strains can only be captured in a crude way. No elasticity and DRX
EVPFFT	Polycrystal	Full-field	Provides effective elastovisco-plastic behavior and local intra-crystalline fields, 2D and 3D	Microstructural effects on local fields distribution and their evolution during transient creep	Limited to small strains, no microstructure evolution yet, no DRX
VPFFT - Elle	Polycrystal	Full-field	Provides local intra-crystalline fields and microstructure evolution in 2D	Couple local field predictions to DRX mechanisms in the secondary creep regime	No account for local field evolution during transient creep. Limited to 2D (Elle). Rough update of the dislocation field during DRX

Table 3: Summary of techniques application domains, interest and limitations. DRX: dynamic recrystallization, VPFFT: FFT-based formulation for Visco-Plastic polycrystals , EVPFFT: FFT-based formulation for Elasto-Visco-Plastic polycrystals

Model	Scale	Type	Applications	Advantages	Limitations
Mixture with Continuous Diversity (MCD)	Large scale	General continuum theory	General overview of the interactions between microstructure evolution, DRX and ice flow	Effects of microstructure and its evolution via internal variables. Secondary and tertiary creep regimes. Thermodynamically consistent	Mathematically complex.  Not implemented numerically yet
CAFFE	Large Scale	Continuum	Provides effective visco-plastic behavior on the large scale, including fabric development	Captures the effects of fabric development in the secondary and tertiary creep regimes, easy to implement	Stress and strain rate are colinear (scalar effective viscosity). Limited to fabrics represented by a multipole expansion up to fourth order only.
GOLF law	Polycrystal	Phenomen- ological orthotropic non-linear law	Provides orthotropic viscous behavior and fabric development	Efficient, easy-to-use and able to reproduce the response of micro-macro models	Orthotropic fabric restricted to the assumed closure approximation. Non-linear case not validated against micro-macro models yet
Elmer/Ice	Large scale	FE code including GOLF and CAFEE law	Flow of anisotropic polar ice and its fabric evolution	Fabric evolution consistent with the stress field and strain-rate field	Can not be used to simulate the localization of the deformation (diffusion of the fabric induced by interpolation)

Table 4: Summary of techniques application domains, interest and limitations. DRX: dynamic recrystallization