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Abstract. The verification and calibration of a new quasi-linear transport model with a large database of gyrokinetic turbulence simulations is presented in this paper. In a previous paper [1], a model for the saturated spectrum of electric potential fluctuations was developed based on the properties of the non-linear 3D spectrum. In this paper, a modification to the overall multiplicative factor in this model is found to be necessary to improve the fit to scans of the temperature and density gradients and safety factor. The error in the fit of the quasi-linear fluxes of electron and ion energy fluxes is significantly better than for previous saturation models. The spectral shift model for the impact of equilibrium $E \times B$ velocity shear [2] and the zonal flow mixing model for electron-scale turbulence [3] are both revised to be compatible with this new model. The models for the loss of bounce averaging and electron collisions in the TGLF reduced linear equations [4] are also changed to improve the linear eigenmodes.

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1. Introduction

In a previous paper [1] a new model for the saturated 3D spectrum of electric potential fluctuations from gyrokinetic turbulence simulations was presented. This model was built, for the first time, using the linear eigenmodes computed with the same gyrokinetic code (CGYRO [5]) as the non-linear simulations. This eliminates the uncertainty of the accuracy of the linear eigenmodes computed with a reduced model like TGLF [4]. The 3D potential fluctuation spectrum was examined, including the poloidal (k_y) , and radial (k_x) wavenumbers, and the dependence on the poloidal angle (θ) . The Miller geometry formalism for shaped toroidal magnetic flux surfaces [6],[7] was used. If was found that the geometric metrics of the perpendicular wave vector, normalized to the physical ion gyroradius, provided the functions needed to fit the poloidal angle dependence of the non-linear fluctuation intensity. These functions capture the flux surface shape dependence well. It will be shown in this paper that this same model, with only an overall re-calibration, fits a much larger set of gyrokinetic simulations.

The database of gyrokinetic simulations is described in Section 2. Fundamental tests of the quasi-linear approximation are presented in Section 3. The model for the saturation of the zonal (axisymmetric) potential fluctuations is tested in Section 4 and found to agree very well with the CGYRO cases. The details of the new model for the finite toroidal modes ($k_y > 0$) are presented in Section 5. The verification and calibration of the saturation model to the CGYRO simulation database is presented in Section 6. In Section 7, improvements to the TGLF linear fluid equation models for electron collisions [8] and loss of bounce averaging [4] are presented. The implementation of the new (SAT2) saturation model in the quasi-linear TGLF transport flux calculation is verified with the CGYRO database in Section 8. The spectral shift model [2] for equilibrium $E \times B$ velocity shear suppression of the turbulence is recalibrated for the SAT2 model in TGLF in Section 9. The zonal flow mixing model for electron scale turbulence [3] is calibrated for TGLF-SAT2 in Section 10. A summary is given in Section 11 followed by an appendix. The appendix discusses the transformation properties of quasi-linear fluxes and the summation measure for the fluxes.

2. Gyrokinetic turbulence simulation database

In order to verify and calibrate the saturation model, a set of CGYRO turbulence simulations, composed of 13 scans about the GA standard (GASTD) case, were collected. The parameters of the GASTD are: $a/L_{n_e} = 1.0$, $a/L_{n_i} = 1.0$, $a/L_{T_e} =$ $3.0, a/L_{T_i} = 3.0, \hat{s} = 1.0, q = 2.0, \kappa = 1.0, r/a = 0.5, R/a = 3.0, \delta = 0.0, \Delta = dR/dr =$ $0.0, a\nu_{ee}/c_s = 0.1, T_i/T_e = 1.0, \beta_e = 0.0005$. There are 64 cases in this database but only 52 are unique. Most scans are for a single parameter but scans 12 and 13 vary magnetic shear and safety factor together such that $\hat{s} = q/2$. All are pure deuterium plasma with

Scan	Parameter	Range	Cases	Fixed
1	κ	1.0, 2.0	1-5	$\nu_{ee}a/c_s = 0.05$
2	Δ	0.0, -0.5	6-10	$\nu_{ee}a/c_s = 0.05$
3	$ u_{ee}$	0.05, 1.0	11 - 17	
4	a/L_T	1.25, 3.5	18-23	
5	a/L_n	0.0, 3.0	24 - 30	
6	$ u_{ee}$	0.5, 1.0	31 - 37	$a/L_n = 3.0$
7	r/a	0.25, 1.5	38-43	
8	R/a	2.0, 3.0	44-46	
9	T_i/T_e	0.5, 1.5	47-49	
10	\hat{s}	0.166, 1.5	50 - 54	
11	q	1.5, 3.0	55 - 57	
12	q	1.6, 4.0	58-61	$\hat{s} = q/2$
13	q	1.6, 3.0	62-64	$\hat{s} = q/2, a/L_n = 3.0$

 Table 1. CGYRO gyrokinetic turbulence simulation database

 Scan
 Parameter
 Range
 Cases
 Fixed

equal density, density gradients and temperature gradients for electrons and ions. Note that scan 7 is physical because the length "a" is arbitrary and $r/R \leq 1.5/3.0$. The first two scans were used to determine the geometry dependence of the 3D saturation model [1]. Many of the rest of the scans were part of a study published by E. Belli [9]. The typical resolution used in these CGYRO turbulence simulations is as follows: 16 toroidal modes, 128 radial modes, 8 energy nodes, 16 pitch angle nodes, 28 poloidal angles $dk_y = 0.067$, BOX_SIZE=6. The multi-species Sugama collision model [10] was used. Departures from the GASTD parameters held fixed for the scan are indicated in the last column of Table 1. All of the CGYRO simulations were run long enough to fully saturate (at least 1200 a/ c_s time units).

3. Fundamental tests of quasi-linear theory

The turbulent energy flux (Q_a) and particle flux (Γ_a) through a radial flux tube for the gyrokinetic simulation are computed for each species (subscript a) by the formulas

$$Q_a = \sum_{k_y} \sum_{k_x} \frac{3}{2} \langle Re[ik_y \frac{ae\phi_{k_x,k_y}^*}{\rho_s T_e} \frac{a\tilde{p}_{a,k_x,k_y}}{\rho_s n_e T_e}] \rangle_{t,\theta} = \sum_{k_y} Q_{ak_y} \tag{1}$$

$$\Gamma_a = \sum_{k_y} \sum_{k_x} \langle Re[ik_y \frac{ae\phi^*_{k_x,k_y}}{\rho_s T_e} \frac{a\tilde{n}_{a,k_x,k_y}}{\rho_s n_e}] \rangle_{t,\theta} = \sum_{k_y} \Gamma_{ak_y}$$
(2)

Here the gyro-Bohm normalizations of the CGYRO code were used [5]: $c_s = \sqrt{T_e/m_D}$, $a = \text{minor radius at separatrix}, \ \rho_s = c_s/\Omega_s, \ \Omega_s = eB_{\text{unit}}/(m_Dc), \ B_{\text{unit}} = \frac{qd\psi}{rdr}$ [7] for poloidal magnetic flux ψ , and m_D = deuterium mass. The gyro-Bohm normalization for energy flux in these units is $n_e T_e c_s (\rho_s/a)^2$. The energy flux is produced by a radial $E \times B$ drift driven by electric potential fluctuations $\tilde{\phi}_{k_x,k_y}$ that is averaged with the fluctuations of the pressure moment of the species distribution function (\tilde{p}_{a,k_x,k_y}) . The angle bracket represents a time average (t) and a flux surface average over the poloidal angle (θ) . There is also a sum over the normalized radial wavenumber (k_x) which is equivalent to a radial average over the periodic flux tube box.

The quasi-linear approximation to the turbulent flux evaluates the non-linear flux formula with the most unstable eigenmode at each poloidal wavenumber $\delta \phi_{k_y}$. Typically the most unstable eigenmode is for $k_x = 0$. The quasi-linear weights of a linear eigenmode are defined by

$$WQ_{ak_y}^{QL} = \frac{3}{2} \frac{\left\langle Re[ik_y \frac{ae\delta\tilde{\phi}_{k_y}^*}{\rho_s T_e} \frac{a\delta\tilde{p}_{a,k_y}}{\rho_s n_e T_e}] \right\rangle_{t,\theta}}{\left\langle \left| \frac{ae\delta\tilde{\phi}_{k_y}}{\rho_s T_e} \right|^2 \right\rangle_{t,\theta}}$$
(3)

$$W\Gamma^{QL}_{ak_y} = \frac{\left\langle Re[ik_y \frac{ae\delta \tilde{\phi}^*_{k_y}}{\rho_s T_e} \frac{a\delta \tilde{n}_{a,k_y}}{\rho_s n_e}] \right\rangle_{t,\theta}}{\left\langle \left| \frac{ae\delta \tilde{\phi}_{k_y}}{\rho_s T_e} \right|^2 \right\rangle_{t,\theta}}$$
(4)

The quasi-linear approximation is valid if the phase angle between potential and velocity moments of the distribution function in the non-linear turbulence preserves the linear eigenmode phase angles contained in the quasi-linear weights. This manifests itself as a preservation of the cross-phase between different velocity moments of the distribution function, like electron density and electron temperature, that have been measured in experiments [11], [12] and found to agree with non-linear and quasi-linear calculations. The phase preservation can also be tested in gyrokinetic turbulence simulations [13], [14], [15], [16]. In this paper the quasi-linear intensity spectrum needed to exactly match the electron energy flux will be computed. This will then be used to compute the ion energy and particle fluxes from their respective quasi-linear weights multiplying this intensity spectrum. The difference between these quasi-linear fluxes and the non-linear fluxes is a measure of the error in the quasilinear approximation originating from the quasi-linear weight. This test will be quantitatively verified for the 54 cases in the database. Note that the small contribution due to the magnetic field fluctuations is not included in the CGYRO fluxes.

Because the non-linear flux contribution for each k_y are summed over k_x and flux surface averaged, it is not possible to directly compute the contribution to the intensity from a single linear eigenmode. Instead, the quasi-linear intensity $I_{k_y}^{QL}$ required to make the quasi-linear flux exactly equal to the non-linear flux will be used. It is computed from the electron energy flux by the formula:

$$I_{k_y}^{QL} = \frac{Q_{ek_y}}{\Delta K_{iy} W Q_{ek_y}^{QL}} \tag{5}$$

The poloidal wavenumber grid interval ΔK_{iy} , defined in Appendix A, is needed in Eq. 5 to make the intensity independent of the k_y grid. If the quasi-linear approximation is



Figure 1. The ion energy flux (left) and particle flux (right) contributions for each k_y from the non-linear CGYRO runs (black) and computed from the quasi-linear weights (gray) times the intensity from the electron energy flux (Eq. 5) for the GASTD case with density gradient $a/L_{n_e} = a/L_{n_i} = 1$ (upper) and 3 (lower)

valid, it should not matter which plasma species (subscript label a = e,i), or transport channel, is used to compute this intensity since the species dependence is accounted for in the quasi-linear weight. This is illustrated in Fig 1 where the flux contribution to each poloidal wavenumber k_y is plotted for CGYRO (black) and the quasi-linear weight multiplied by $\Delta k_y I_{k_y}^{QL}$ (gray) for ion energy flux (left) and particle flux (right). The electron energy flux match is exact by construction. For the GASTD case (top) the match for ion energy flux is good but the particle flux peak is lower for the quasi-linear flux indicating that the non-linear particle flux departs somewhat from the phase of the quasi-linear particle flux weight for this case. For a higher density gradient (bottom) the match is very good for both ion energy (left) and particle flux (right) in Fig. 1. Note that the particle flux is much higher for higher density gradient (lower right) than the lower density gradient (upper right). It is consistently found for the CGYRO database that the quasi-linear weight for particle flux is not as well preserved for low particle flux cases.

In order to be able to compare the error due to the quasi-linear weights with the errors reported below for the full quasi-linear flux model, the fractional error from using the quasi-linear intensity computed from the electron energy flux (Eq. 5) to compute the ion energy flux and particle fluxes for the database cases is computed in the following



Figure 2. The error in the QL weight computed from the ion energy flux (left) and particle flux (right) for the 13 scans of the database Table 1

way:

$$\sigma_{WQ_i} = \sqrt{\frac{\sum_m \left(Q_i - \sum_{k_y} \Delta K_{iy} I_{k_y}^{QL} W Q_{ik_y}^{QL}\right)^2}{\sum_m Q_i^2}} \tag{6}$$

$$\sigma_{W\Gamma_e} = \sqrt{\frac{\sum_m \left(\Gamma_e - \sum_{k_y} \Delta K_{iy} I_{k_y}^{QL} W \Gamma_{ek_y}^{QL}\right)^2}{\sum_m \Gamma_e^2}}$$
(7)

The summation index "m" labels the members of each scan in the database (column 3 of Table 1).

In Fig. 2 are shown the errors for the ion energy flux (left) and particle flux (right) for scans in the database. The root mean squared average error for the whole database is 5.4% for the ion energy flux and 19.8% for the particle flux. The high density gradient scans (6 and 13) have the lowest particle flux errors. These errors from the quasi-linear weights set a floor for the quasi-linear flux model since the model for the saturated intensity will have its own errors added in. The degree of error from non-linear departures from the quasi-linear weights is quite low for energy fluxes and particle fluxes of sufficient size.

4. Saturation of the zonal potential fluctuations

The time dependent zonal $(k_y = 0)$ electric potential fluctuations play a strong role in the saturation of the gyrokinetic turbulence. Study of multi-scale (electron+ion gyroradius scales) turbulence simulations [17], [18] demonstrated that $E \times B$ flow shear due to the zonal potential could not compete with the linear growth rate at electron scales $(k_y > 1)$. A new paradigm for saturation through zonal flow mixing [3] was proposed. The zonal flows mix the most unstable modes $(k_x = 0)$ with subdominant, or stable modes, across the k_x spectrum, by coupling modes with the same poloidal wavenumber k_y to modes with different radial wavenumbers k_x . This mixing can compete with the linear growth rate at all scales since the zonal flow mixing rate $V_{ZF}k_y$



Figure 3. The RMS zonal velocity computed from the CGYRO simulations vs the maximum of the linear growth rate divided by the poloidal wavenumber

scales with the poloidal wavenumber of the mixed mode. A simple detailed balance between the linear growth rate and the zonal flow mixing rate provides an estimate of the saturated amplitude of the RMS zonal flow velocity.

$$V_{ZF} = \max[\gamma_{k_y}/k_y] \tag{8}$$

Here, the linear growth rate γ_{k_y} (normalized by c_s/a) is for the most unstable mode at each k_y . The RMS zonal flow velocity can be computed directly from the non-linear saturated spectrum

$$V_{ZF} = 0.5 \sqrt{\sum_{k_x} k_x^2 \left| \frac{a e \tilde{\phi}_{k_x,0}}{\rho_s T_e} \right|^2} \tag{9}$$

The RMS zonal flow velocity V_{ZF} is independent of k_y so its saturated magnitude is set by the maximum drive in Eq. 8. This saturation rule for V_{ZF} (Eq. 8) is shown to be well satisfied for the database in Fig. 3 The line in Fig. 3 is a fit to the data giving a linear coefficient of 1.12. The coefficient of (0.5) in Eq. 9 would thus be a better fit with 0.45. The saturation model below for the finite k_y spectrum does not depend on this coefficient. The same data as in Fig. 3 is plotted for all of the database cases in Fig 4. The lowest values are for the temperature gradient scan 4. The highest values are for scan 7 (r/a). A failure of the zonal flow saturation rule (Eq. 9), for a particular gyrokinetic simulation, would be an indication that zonal flow mixing is not the dominant saturation mechanism or that the simulation has not yet reached saturation.



Figure 4. The RMS zonal velocity computed from the CGYRO simulations vs the maximum of the linear growth rate divided by the poloidal wavenumber

In the previous study of the 3D spectrum [1] it was observed that the RMS width of the radial wavenumber spectrum for finite k_y was limited to be larger than the width of the zonal potential spectrum. An accurate method for fitting the radial wavenumber spectrum with a Lorentzian model was employed to determine this width. The error in this fit tends to be larger for the zonal potential than the rest of the spectrum because the zonal potential spectrum sometimes narrows at low k_x [1]. For the first two scans, that were previously studied, the width of the zonal potential spectrum tracked the poloidal wavenumber at the maximum zonal flow drive $k_{xZF} \approx k_{ymax}/|\nabla r|_0$ where the geometric factor is evaluated at the outboard midplane. For the wider database of this paper, it is found that this approximate model needs to be generalized to:

$$k_{xZF}^{model} = 0.83\sqrt{(R/Lp12)/(q/2))}k_{ymax}/|\nabla r|_0$$
(10)

Here R is the major radius at the center of the flux surface, $R/L_p = -(R/P)dP/dr$ where P is the total pressure, and q is the safety factor. This model is shown for all of the cases along with the spectral width k_{xZF} computed from the saturated zonal potential spectrum in Fig. 5. The lowest width k_{xZF} is for the temperature gradient scan 4 (cases 18-23). The highest widths are for high density gradient and high collision frequency (cases 36-37). The fit of the model is not tight but it is shown here because it motivates the factors needed to calibrate the overall fluxes to the CGYRO database that will be presented in the next section.



Figure 5. The RMS width of the zonal potential spectrum and the model Eq. 10

5. The new saturation model

The model for the saturated quasi-linear intensity developed in Ref. [1] is

$$I_{k_y}^{\text{model}} = \left\langle G^2(\theta) \right\rangle_{\theta} \left(\frac{\gamma_{k_y}^{\text{model}}}{k_x^{\text{model}} k_y} \right)^2 \tag{11}$$

All of the θ dependence is absorbed into $G(\theta)$ that was fit to the theta dependence of the square amplitude of the peak potential fluctuation spectrum at $k_x = 0$ [1].

$$G^{2}(\theta) = d_{1}G_{1}(\theta) \text{ for } k_{y} < k_{ycut}$$

= $\left(d_{1}G_{1}(\theta)k_{ycut} + b_{3}d_{2}G_{2}(\theta)(k_{y} - k_{ycut})\right)/k_{y} \text{ for } k_{y} \ge k_{ycut}$ (12)

where $b_3 = 2.4$ and the coefficients d_1, d_2 are

$$d_1 = \left(\frac{B_{T_0}}{B(0)}\right)^4 \tag{13}$$

$$d_2 = 1/G_q(0)^2 \tag{14}$$

here $B(0)|_{r=0} = B_{T_0}$ is the toroidal magnetic field at the flux surface center and $G_q(\theta) = |\nabla r| B_{\text{unit}}/B(\theta)$. Two geometric shape functions are:

$$G_1 = \left(\frac{B(0)}{B(\theta)}\right)^4 \tag{15}$$

$$G_2 = \left(\frac{G_q(0)}{G_q(\theta)}\right)^4 \tag{16}$$

The cutoff is modeled by $k_{ycut} = b_0 k_{ymax}$ where k_{ymax} is the value of k_y where γ_{k_y}/k_y is maximum and the best fit parameter was determined to be $b_0 = 0.72$.

The model (k_x^{model}) for the width of the potential radial wavenumber spectrum at finite k_y was fit to the k_x width of the turbulence spectrum at the outboard midplane $\theta = 0$. The model is given by:

$$k_x^{\text{model}} = \frac{B_{\text{unit}}}{B_{\text{norm}}} k_{\text{ycut}} / |\nabla r|_0 \quad \text{for} \quad k_y < k_{\text{ycut}}$$

$$= \frac{B_{\text{unit}}}{B_{\text{norm}}} \left(k_{\text{ycut}} / |\nabla r|_0 + b_1 (k_y - k_{\text{ycut}}) G_q(0) \right) \quad \text{for} \quad k_y \ge k_{\text{ycut}}$$
(17)

Here B_{norm} is the arbitrary external magnetic field normalization which is chosen to be B_{unit} for CGYRO. The best fitting coefficient was determined to be $b_1 = 1.22$. This model differs from the previous saturation model (SAT1) [3] in that the k_x width does not continue to scale with k_y below k_{ycut} but rather flattens out. This is conjectured to be the effect of the finite width of the zonal potential spectrum. There can be more complex structure to the k_x width in the region $k_y \leq k_{ymax}$ that is not captured completely by this model. The dependence of the slope of k_x^{model} for $k_y \geq k_{ycut}$ on the geometric factor G_q is also new to this SAT2 model.

The effective non-linear mixing rate $\gamma_{k_y}^{\text{model}}$ in Eq. 11 is modeled by

$$\gamma_{k_y}^{\text{model}} = b_2 \gamma_{k_y} \text{ for } k_y < k_{\text{ymax}}$$
$$= b_2 \gamma_{max} \text{ for } k_y \ge k_{\text{ymax}}$$
(18)

Here γ_{max} is the value of the linear growth rate at k_{ymax} where the peak in γ_{k_y}/k_y occurs. The offset for the total energy fluxes summed over scans 1 and 2 in Ref. [1] is set to zero for $b_2^2 = 7.94$. This model differs from the SAT1 model [3] in that it is simply b_2 times the linear growth rate for $k_y \leq k_{ymax}$. The flattening of the k_x width model suppresses the intensity at low k_y making the zonal flow mixing subtraction included in SAT1 for $k_y \leq k_{ymax}$ unnecessary.

The reader is referred to Ref. [1] for the procedure used to determine this model from the 3D potential spectrum properties. The four fitting coefficients (b_0, b_1, b_2, b_3) were adjusted to minimize the absolute RMS difference between the CGYRO and QL energy flux contributions at each k_y for the first two scans in the database. These coefficients were not changed other than the overall renormalization given in Eq. 19. In order to fit the wider database of this paper, it was found that the overall coefficient needed to be modified to

$$b_2^2 \to 8.44 \left(\frac{q}{2} \frac{12L_p}{R}\right) \tag{19}$$

The value 8.44 is determined by forcing the offset in the sum of electron and ion energy fluxes to be zero for the whole database. This modification may be related to the trend for the width of the zonal potential spectrum found in Eq. 10 since $k_x^{\text{model}} \approx k_{xZF}$ at low k_y . Note that the GASTD case $(q = 2, R/L_p = 12)$ is taken as a reference point.



Figure 6. The fractional error for the 13 scans in the CGYRO database. Ion energy flux (top), electron energy flux (middle) and particle flux (bottom)

6. Verification and calibration of the quasi-linear fluxes

The quasi-linear model fluxes are computed from the formulas:

$$Q_a^{QL} = \Delta K_{iy} I_{ky}^{\text{model}} W Q_{aky}^{QL} \tag{20}$$

$$\Gamma_a^{QL} = \Delta K_{iy} I_{k_y}^{\text{model}} W \Gamma_{ak_y}^{QL} \tag{21}$$

The linear growth rate and quasi-linear weight spectrum are computed using CGYRO. The fractional error for these quasi-linear fluxes compared to the CGYRO fluxes is computed for each scan by the formula:

$$\sigma_{Q_a^{QL}} = \sqrt{\frac{\sum_m \left(Q_a - Q_a^{QL}\right)^2}{\sum_m Q_a^2}} \tag{22}$$

$$\sigma_{\Gamma_a^{QL}} = \sqrt{\frac{\sum_m \left(\Gamma_a - \Gamma_a^{QL}\right)^2}{\sum_m \Gamma_a^2}}$$
(23)

The fractional errors for all of the database are given in Fig. 6. The RMS average of these errors is 12.0% for Q_i , 11.7% for Q_e and 20.8% for Γ_e . These are almost a factor of 2 lower errors than previous saturation models and are not far from the uncertainty of the gyrokinetic simulations due to finite time averaging and grid resolution. The particle flux error due to the quasilinear weights (19.8%) is the primary contribution for this channel. The quasi-linear weight error is about half of the total error for energy fluxes.

In Fig. 7 are plotted the ion energy flux (top) electron energy flux (middle) and particle flux (bottom) for CGYRO (black) and the QL model (gray) for all of the 64 cases in the database. The cases are numbered sequentially in the order they appear in the database Table 1.

The modification of the overall fluxes with pressure gradient was needed for the scans in temperature gradient (4) and density gradient (5). These scans could also be fit by modifying the overall coefficient to vary with the ratio of γ_{max} to a proxy for the geodesic acoustic mode frequency v_{thi}/R such that the intensity scaled like the first power of γ_{max} instead of the second power. However, this choice spoils the good agreement with other scans (e.g. collisions, magnetic shear). The pressure gradient factor essentially changes the flux surface label from "r" to "P" in the linear gradient drift terms. Since the total pressure gradient is fundamental to the radial force balance of the plasma, it defines a spacing between flux surfaces that perhaps has an impact on the turbulence. If this is an equilibrium effect, then the total pressure including fast ions should be used. This needs to be verified. The pressure gradient factor is introduced to improve the fit for these 54 cases but a larger database may invalidate it. The improvement of the fit by introducing an overall factor of the safety factor is not surprising, given the empirical scaling law trend that energy confinement improves with current rather than magnetic field. Its origin may be due to the zonal potential width but the physics of this is not understood. Fortunately, the more complicated fit for the effective mixing rate (Eq. 18) and radial wavenumber spectral width (Eq. 17) did not require revision from the fit in Ref. [1].

7. Improvements to the TGLF linear eigensolver

The use of the linear eigenmodes from CGYRO allowed the fitting of a saturation model to the non-linear CGYRO simulations without concern for the accuracy of the



Figure 7. The fluxes for CGYRO (black) and the QLSAT2 model (gray) for all of the cases. Ion energy flux (top), electron energy flux (middle) and particle flux (bottom)

linear eigenmodes. The problem with this approach is that the calculation of the linear CGYRO eigenmodes is very slow compared to a reduced gyro-fluid model like TGLF [4]. The CGYRO linear calculation is also an initial value run so it only finds the most unstable mode, whereas TGLF is a matrix eigensolver that finds all of the unstable branches of its equations. Including subdominant modes is essential for helping the flux smoothly cross branch jumps. Initial value runs have convergence difficulty near the linear stability threshold or when there are two modes with similar growth rates but different frequencies. These issues make it impractical to use CGYRO (or any initial value gyrokinetic code) as the linear eigenmode solver in a transport code.



Figure 8. The CGYRO (black) and TGLF for the X2 (dashed) and X3 (gray) collision models linear growth rate/ k_y (left) and frequency/ k_y (right) for the GASTD case with $a/L_{n_e} = a/L_{n_i} = 3.0$

The linear eigenmodes of TGLF were found to lack accuracy for the high density gradient cases in the database. For example, the spectrum of the growth rate (left) and frequency (right), divided by the poloidal wavenumber k_y for the GASTD case parameters but with higher density gradients $(a/L_{n_e} = a/L_{n_i} = 3.0)$ is shown In Fig. 8. The black curves are the calculation using CGYRO with the Sugama multi-species collision operator. The Dashed curves are for TGLF with the collision model of Ref. [8] (denoted X2 here for option XNU_MODEL=2). The TGLF equations have bounce averaged trapped particles for all species. The X2 collision model is only pitch angle scattering of electrons. The dominant term in the TGLF collision model is from the gradient of the distribution function at the trapped-passing boundary which is modeled by ($\xi = v_{\parallel}/v$)

$$\nu_{ee} \frac{df_e}{d\xi}\Big|_{\xi=\xi_t} = c_b \nu_{ee}^{(1-\sigma_b)} (v_{the}k_{||})^{\sigma_b} \tilde{f}_e \tag{24}$$

Where $\xi_t = \sqrt{1 - B_{\min}/B_{\max}}$ is the trapped-passing boundary value. For the X2 model $c_b = 0.114$ and $\sigma_b = 0.5$. The linear growth rate in Fig. 8 for the X2 model (dashed) is too large across most of the spectrum. This 21% larger peak in γ_{k_y}/k_y translates to about 47% larger fluxes for TGLF with the X2 model compared to CGYRO using the saturation model of this paper. Increasing the coefficient c_b of the trapped-passing boundary collision term lowers the linear growth rate at low k_y but this shifts the peak to higher k_y and does not help lower the higher k_y growth rate.

The TGLF equations have a model for the loss of bounce averaging of the Landau resonance by trapped particles [4]. The model attempts to estimate how far into the trapped region of velocity space the particles can experience Landau damping and hence should be part of the circulating particle fluid. Assuming that the Landau resonance condition determines the frequency gives the relation

$$\omega = k_{||}v_{||} = k_{||}Rqd\theta/dt \tag{25}$$

Integrating this equation in time gives $\omega \Delta t = k_{\parallel} Rq \Delta \theta$. A trapped particle keeps the same sign of its velocity while traveling between bounce angles θ_B . A minimum

requirement for a trapped particle to be able to average the Landau resonance is that it can change the sign of its velocity before the parallel electric field of the wave changes sign. Hence it must travel more than half an orbit $(2 \theta_B)$ in half a wave period $(\omega \Delta t = \pi)$. This gives the condition for a trapped particle to be able to Landau average

$$\Delta \theta = \frac{\pi}{Rqk_{||}} > 2\theta_B \tag{26}$$

This gives a condition for the maximum bounce angle for trapped particles that can average the Landau resonance θ_{LA} [4]

$$\theta_{LA} = MIN\left(\frac{\pi}{2Rqk_{||}},\pi\right) \tag{27}$$

The effective boundary for trapped particles that can bounce average the Landau resonance is thus $\xi_{LA} = \sqrt{1 - B_{\min}/B(\theta_{LA})}$. In practice this condition is used in TGLF to reduce the trapped fraction using an estimate of the parallel wavenumber k_{\parallel} . The eigenvalue solution procedure for TGLF adjust the parallel wavenumber with just a Gaussian wavefunction with the first 2 Hermite polynomials to find the maximum linear growth rate. The wavefunction is then refined with more Hermite polynomials (typically 4). For high magnetic shear, the parallel wavenumber found is larger and the effective trapped particle boundary is reduced. This tracks the linear growth rate for trapped electron modes well [4]. The TGLF problem of too high a linear growth rate at high density gradient is conjectured to be due to the higher density gradient drift frequency $(k_y a/L_{n_e})$ causing a loss of bounce averaging. In order to include this effect, the above model is modified in the following way

$$\theta_{aLA} = MIN\left(\frac{\pi}{2Rqk_{a||}^{\text{eff}}}, \pi\right) \tag{28}$$

$$k_{a||}^{\text{eff}} = |k_{||}| + 3.0 \frac{B_{\min}}{B_{\max}} \left| \frac{ak_y}{v_{th_a} L_{n_a}} \right|$$
(29)

The coefficient 3.0 in Eq. 29 was determined by fitting to the CGYRO linear growth rate spectrum at higher density gradient. The new model for the loss of Landau averaging (Eq. 29) is species dependent, since the diamagnetic drift frequency $k_y a/l_{n_a}$, and the thermal velocity $v_{tha} = \sqrt{(T_a/m_a)}$, are species dependent. The trapped ions lose bounce averaging at a lower k_y than the electrons in this model, which makes a difference at the lowest k_y range. The trapped-passing boundary collision model (Eq. 24) was adjusted to $c_b = 0.315$, $\sigma_b = 0.34$ in order to match the decay of the CGYRO growth rate spectrum with collisions. This change in the electron collisions and the model for the loss of bounce averaging is the X3 model (option XNU_MODEL=3). The gray curves in Fig. 8 show the improvement in the TGLF fit for the X3 model. It was found that this model was good up to about $\nu_{ee}a/c_s = 1.0$. Above this value for the electron collision frequency, the low k_y modes are over-stabilized. A new eigensolver for TGLF that does not employ the bounce averaging approximation is under development. This will eliminate the need for these ad-hoc adjustments to bounce averaging.



Figure 9. The error between the TGLF fluxes and CGYRO for each scan. Ion energy flux (top), electron energy flux (middle) and particle flux (bottom)

8. TGLF implementation of the QL model

With the X3 collision model, and the new SAT2 saturation model presented in this paper, TGLF provides a good quasi-linear model of the CGYRO non-linear fluxes in the database. The error between the TGLF fluxes and CGYRO for the scans are shown in Fig. 9. The RMS average error over all of the scans is: Qi 15.6%, Qe 14.2% ,and Γ_e 35.6%. This is not quite as good as found using the CGYRO eigenvalues but is significantly better than previous saturation models. The number of Hermite polynomials for the ballooning mode wavefunction in TGLF was increased from 4 to



Figure 10. The fluxes for CGYRO (black) and the TGLF model (gray) for all of the cases. Ion energy flux (top), electron energy flux (middle) and particle flux (bottom)

6 for the four cases with $q \ge 3$ (cases 57, 60, 61, 64). This reduced the overshoot of the growth rates for these cases. This inaccuracy of the TGLF eigenmodes with the standard 4 Hermite basis functions may be why previous saturation models that were fit using TGLF eigenmodes did not appear to require an overall multiplication by the safety factor as was found for the present model. In Fig. 10 is shown the ion energy (top), electron energy (middle) and particle (bottom) fluxes for all of the cases. Compared to the QL fluxes computed with GYRO eigenmodes in Fig. 7 the TGLF particle flux is reduced, especially for the scan 7 (cases 38-43) that scans the local radius r/a and hence scans the trapped fraction. This suggest a deficit in the trapped electron contribution to the particle flux for TGLF.

9. Calibration of the spectral shift model for ExB velocity shear

The spectral shift model for the effect of equilibrium $E \times B$ velocity shear $(\gamma_{E \times B} = (q/r)d(d\phi/d\psi)/dr)$ [7] on gyrokinetic turbulence [2] captures the shift in the peak of the 2D flux surface averaged potential spectrum away from $k_x = 0$ to a finite value k_{x0} . The direction and size of this shift is determined by the $E \times B$ velocity shear. This spectral shift produces an ion stress contribution to momentum transport by breaking the parity of the eigenmode [19]. However, this shift in the eigenmode does not reduce the linear growth rate enough to account for the reduction in the fluxes. The peak of the potential spectrum is reduced by two effects. One is the change in the radial correlation length that was previously identified as a suppression mechanism [20], [21]. This is manifest in the spectral shift model as a change in the width of the radial wavenumber spectrum and a tilt in the radial orientation of the turbulent eddies [19]. This effect alone is too weak to account for the reduction in fluxes with $E \times B$ velocity shear in the turbulence simulations [22]. The second effect is a temporal reduction in the effective mixing rate $\gamma_{k_y}^{\text{eff}}$. Both of these reductions in the peak of the potential were shown [2] to be directly related to the shift k_{x0} in the spectrum by the formula :

$$\Phi(k_x, k_y) = \frac{\Phi_0(0, k_y)}{\left(1 + \left(\frac{k_{x0}}{k_x^{\text{model}}}\right)^2 + \left(\frac{(k_x - k_{x0})}{k_x^{\text{model}}}\right)^2\right) \left(1 + (\alpha_x \frac{k_{x0}}{k_x^{\text{model}}})^{\sigma_x}\right)}$$
(30)

Here Φ is the flux surface and time averaged gyro-Bohm normalized potential fluctuation amplitude and $\Phi_0(0, k_y)$ is the peak of the potential in the absence of equilibrium $E \times B$ velocity shear. The left hand factor in the denominator in Eq. 30 accounts for the broadening of the radial wavenumber spectral width by the spectral shift k_{x0}

$$k_x^{\text{net}} = \sqrt{\frac{\sum_{k_x} (k_x - k_{x0})^2 \Phi^2}{\sum_{k_x} \Phi^2}} = k_x^{\text{model}} \sqrt{1 + \left(\frac{k_{x0}}{k_x^{\text{model}}}\right)^2}$$
(31)

The radial wavenumber spectral width in the absence of equilibrium $E \times B$ velocity shear is given by the model k_x^{model} in Eq. 17. The additional temporal reduction in the peak amplitude is given by the right hand factor in the denominator of Eq. 30.

In the original spectral shift model, the formula for the spectral shift k_{x0} that was determined to fit GYRO simulations has a non-linear dependence on the $E \times B$ velocity shear that also depends on the flux surface shape [19]. Fitting to CGYRO simulations with different flux surface shape was carried out and a linear relation was found to be sufficient. This difference is thought to be due to the fact that CGYRO is spectral in k_x so it has higher accuracy for the gyro-averaging operator (Bessel functions) than GYRO which uses a radial grid. It is possible to use a full Fourier transform to obtain spectral accuracy in GYRO but this was not done in the original simulations used to calibrate the spectral shift model due to the computational expense.



Figure 11. Spectrum of the radial wavenumber shift k_{x0} , computed from CGYRO simulations Eq. 33 (black), and the model Eq. 32 (gray), for a range of values of the equilibrium ExB velocity shear $\gamma_{E\times B}$, for the GASTD case.



Figure 12. The RMS zonal velocity vs $MAX\left[\gamma_{k_y}^{\text{net}}/k_y\right]$ (Eq. 34) for cases: (1) GASTD $\gamma_{E\times B} = 0, 0.1, 0.2, 0.3, 0.4, 0.5, (2) \ \kappa = 1.25, \ \gamma_{E\times B} = 0, 0.1, 0.2, 0.3, (3) \ \kappa = 1.5, \ \gamma_{E\times B} = 0, 0.1, 0.2, 0.3, (4) \ \kappa = 1.75, \ \gamma_{E\times B} = 0, 0.1, 0.2, 0.3, (5) \ \Delta = -0.1, \ \gamma_{E\times B} = 0, 0.1, 0.2, 0.3, (6) \ \Delta = -0.2, \ \gamma_{E\times B} = 0, 0.1, 0.2, 0.3, (7) \ \Delta = -0.4, \ \gamma_{E\times B} = 0, 0.1, 0.2, 0.3$



Figure 13. Scan in the equilibrium ExB velocity shear for the GASTD case. The fluxes for CGYRO (black) and the TGLF model (gray). Ion energy flux (top left), electron energy flux (top right), particle flux (bottom left) and Ion toridal stress (bottom right)

Implementing the spectral shift model in TGLF requires two passes. The first pass computes the linear eigenvalue spectra without $E \times B$ velocity shear. The second pass uses these bare linear growth rates to compute the spectral shift k_{x0}^{model} (Eq. 32). This spectral shift is then used to recompute the linear shifted ballooning eigenmodes to obtain the quasi-linear weights. The shifted ballooning modes produce a small change in the energy and particle weights but produce all of the phase shift required for the ion stress. In the original spectral shift model, the bare linear growth rate at each k_y was used in the formula for the spectral shift at each k_y . This implementation is still used for the older saturation models (SAT_RULE= 0 or 1).

For the new saturation model of this paper (SAT_RULE = 2) the bare growth rates are used to compute the bare values of k_{ymax} and γ_{max} . The spectral shift spectrum is then computed from the formula:

$$k_{x0}^{\text{model}} = 0.32k_y \left(\frac{k_{ymax}}{k_y}\right)^{0.7} \left(\frac{\gamma_{E \times B}}{\gamma_{max}}\right)$$
(32)

This is a reasonable fit to the spectral shift k_{x0} computed directly from the CGYRO



Figure 14. Scan in the equilibrium ExB velocity shear for Elongation (KAPPA) 1.5. The fluxes for CGYRO (black) and the TGLF model (gray) for the GASTD case. Ion energy flux (top left), electron energy flux (top right), particle flux (bottom left) and Ion toridal stress (bottom right)

spectra

$$k_{x0} = \frac{\sum_{k_x} |\tilde{\phi}_{k_y,k_x}|^2 k_x}{\sum_{k_x} |\tilde{\phi}_{k_y,k_x}|^2}$$
(33)

as shown in Fig. 11 for the GASTD case.

The model Eq. 32 is able to fit the spectral shift for a scan in elongation and Shafranov shift without explicit geometry factors. This is a simplification compared to the original model that used the linear bare growth rate to normalize the ExB velocity shear and required a geometric correction.

Next, the bare linear growth rate without $E \times B$ shear is modified by the temporal reduction factor

$$\gamma_{k_y}^{\text{net}} = \frac{\gamma_{k_y}}{\left(1 + \left(\alpha_x \frac{k_{x0}}{k_x^{\text{model}}}\right)^{\sigma_x}\right)} \tag{34}$$

This is then used to recompute the maximum of $\gamma_{k_y}^{\text{net}}/k_y$ to find the net γ_{max} and k_{ymax} that is used in the saturation model (Eq. 30). The net spectral width (Eq. 31) is also used in place of the bare spectral width in the saturation formula.

The zonal potential spectrum does not shift but remains symmetric about $k_x = 0$



Figure 15. Scan in the equilibrium ExB velocity shear for Shafranov shift (SHIFT) -0.2. The fluxes for CGYRO (black) and the TGLF model (gray) for the GASTD case. Ion energy flux (top left), electron energy flux (top right), particle flux (bottom left) and Ion toridal stress (bottom right)

as it must due to the reality condition on the Fourier amplitudes. The RMS average zonal velocity (Eq. 9) is observed to be reduced by the equilibrium $E \times B$ velocity shear. The zonal flow saturation rule remains approximately true if the net growth rate is used as shown in Fig. 12 for a set of 30 cases. The temporal reduction factor was computed using the value of the spectral shift spectrum k_{x0} (Eq. 33) computed from the nonlinear spectrum and the model Eq.17 for the width of the spectrum. This gives about the right level of reduction in zonal flow velocity even though the coefficients $\alpha_x = 1.55$ and $\sigma_x = 2$ were fit to the fluxes not the zonal velocity. This demonstrates that the saturated zonal mixing rate remains in balance with the net effective linear drive as the turbulence is suppressed by the equilibrium $E \times B$ velocity shear.

A novel method for including equilibrium ExB velocity shear in a periodic flux tube was used for these simulations [23]. The equilibrium $E \times B$ velocity is introduced as a sawtooth waveform over the radial periodic box. The magnitude of the flow shear is the same everywhere but the sign flips at the peaks of the sawtooth pattern. This sign flip has no effect on the energy and particle fluxes but reverses the sign of the toroidal stress. The toroidal stress is computed in the central zone of the flux tube where the ExB velocity shear has one sign. The shift in the radial wavenumber spectrum is also computed in this same zone. A good agreement between the k_x spectral code CGYRO using this method and GYRO using a radial grid was shown in Ref. [23] for the GASTD case with adiabatic electrons. The CGYRO results in Fig. 13 are in good agreement with the GYRO results of Ref. [2] for the GASTD case with kinetic electrons but the reduction in the energy fluxes is greater at higher shear for GYRO than CGYRO.

Using the model for the spectral shift (Eq. 32) a good fit to the energy and particle fluxes is obtained for parameters $\alpha_x = 1.55$, $\sigma_x = 2$. Note that this is significantly different than the fit to GYRO simulations [2] with $\alpha_x = 1.15$, $\sigma_x = 4$. This new CGYRO fit is only used for the new saturation model (SAT_RULE = 2) in TGLF.

The shift in the ballooning mode radial wavenumber, that produces the stress, was found to require a multiplier $G_x = -\text{sign}(B_T)0.7/|\nabla r|_0^2$ times k_{x0}^{model} in order to better track the Shafranov shift dependence of the toroidal stress. Note that TGLF and CGYRO have opposite conventions for the direction of the toroidal angle. The CGYRO (black) and TGLF (gray) ion energy flux (top left), electron energy flux (top right), particle flux (bottom left) and ion toroidal stress (bottom right) scans in the equilibrium $E \times B$ velocity shear are shown for the GASTD case in Fig. 13, and cases with elongation $\kappa = 1.5$ in Fig. 14 or with Shafranov shift $\Delta = -0.2$ in Fig. 15.

10. Calibration of the multi-scale model for electron scale energy transport

The final step in the calibration of TGLF with the new saturation model is to extend the model to electron scales. The multi-scale simulations of a C-MOD L-mode plasma by Howard [17] with the GYRO code were used to formulate the first multi-scale saturation model [3] (SAT_RULE = 1). The ion-scale sector of this model is replaced by the new model of this paper. The electron scale modes were found to be suppressed by the zonal flow mixing limiting the effective non-linear mixing rate to the ion scale γ_{max} for $k_y \ge k_{ymax}$ as in Eq. 18. However, if the electron scale peak of γ_{k_y}/k_y is larger than the ion scale peak, the electron scale contribution to the potential can increase yielding a larger contribution to the electron energy flux for $k_y \geq 1$. This effect is visible in Fig. 16 where the GYRO (black) and TGLF-SAT2 (gray) electron energy flux (squares), ion energy flux (circles) and the contribution to the electron energy flux from electron scales $(k_y \ge 1)$ (triangles) are shown for a scans in the ion temperature gradient (left) and electron temperature gradient (right) about an L-mode discharge on the C-MOD tokamak [17]. The increase in the electron scale energy flux as the ion temperature gradient is lowered (left panel) is due to the zonal flow mixing rate $(V_{ZF}k_y)$ lowering below the linear growth rate at electron scales. This effect is built into the saturation model for $k_y \ge k_{ymax}$ by changing the formula Eq. 18 to:

$$\gamma_{k_y}^{\text{eff}} = b_2(\gamma_{max} + MAX[\gamma_{k_y} - \alpha_{ZF}k_yV_{ZF}, 0.0]) \tag{35}$$

The TGLF fluxes with the new saturation model use this formula with $\alpha_{ZF} = 1.1$. The electron-scale saturation model includes a Lorentz distribution weighted average



Figure 16. Scan in the equilibrium ion (Left) and electron (right) temperature gradients for multi-scale GYRO simulations of a C-MOD L-mode [17]. The energy fluxes for GYRO (black) and the TGLF model (gray). Ion energy flux (circle), electron energy flux (square) and electron-scale electron energy flux (triangles)

over k_y of $\gamma_{k_y}^{\text{eff}}$ [3]. This represents the effect of the mixing between modes with different k_y . In the original model, a change in the scaling of the width of the radial wavenumber spectrum for $k_y \geq 1.0$ was included. The new model accounts for this anisotropy of the spectral width through the geometric factors in Eq. 12 and Eq. 17. The original multi-scale GYRO runs output flux surface averaged potential fluctuation spectra so the poloidal angle dependence is unknown. New multi-scale simulations are needed to verify that the poloidal angle dependence of the new model is correct at electron scales. The agreement of the TGLF fluxes with GYRO in Fig. 16 is very good. These simulations include an impurity ion, shaped flux surfaces and a low amount of equilibrium $E \times B$ velocity shear [17],[18].

11. Summary

A new, more accurate, quasi-linear saturation model (SAT2) has been verified and calibrated to a database of CGYRO turbulence simulations in this paper. The model began with exploring the 3D structure of the saturated potential fluctuation spectrum in a previous paper [1]. The SAT2 model was constructed using the linear growth rate and quasi-linear weights from the spectrum of the most unstable modes computed with the same code as the non-linear turbulence simulations. The saturation model is fit to the poloidal wavenumber spectrum of the intensity required to make the quasi-linear approximation to the electron energy flux exact (QL intensity). It was demonstrated that the quasi-linear weights accurately predict the ratio of the ion energy and particle fluxes to the electron energy flux. This quantitative test showed that quasi-linear weights contribute 5.4% of the error in the quasi-linear ion energy fluxes averaged over the whole database of 13 CGYRO scans. The quasi-linear weights contribute 19.8% error to the particle flux due to lower quasi-linear fluxes at low density gradient than from the turbulence. The SAT2 model for the quasi-linear intensity introduces additional error

to the quasi-linear fluxes. The fractional error in the quasi-linear fluxes for the 13 scans is 12.0% for ion energy, 11.7% for electron energy, and 20.8% for particle flux. These errors are lower than previous saturation models fit to gyrokinetic simulations [24], [3].

The zonal flow mixing rule (Eq. 8), for the saturation of the zonal potential fluctuations, was shown to be quite accurate for the whole database. It was even found to work when there is equilibrium $E \times B$ velocity shear if the linear growthrate spectrum is modified by the spectral shift (Eq. 34). This is a central part of the SAT2 model (Eq. 11) spectral shape and peak intensity. The width of the radial wavenumber spectrum, that is conjectured to be determined by the zonal potential mediated coupling of different radial wavenumbers for the same poloidal wavenumber, is the other central element of the SAT2 model. It was found in this larger database that the width of the zonal potential spectrum (Eq. 10) is more variable than previously observed. This deserves further study to understand the physical mechanisms at work. The observed broadening of the radial wavenumber spectral width at low k_y causes a reduction in the intensity of the turbulence. Including this effect in the SAT2 model allowed for a simplification of the model compared to its predecessor [3] while giving a better fit to the quasi-linear intensity at low poloidal wavenumber.

The saturation model is an independent component of a quasi-linear approximation to transport fluxes that can be used with any reduced model for gyrokinetic linear stability to compute quasi-linear fluxes. In this paper, the SAT2 model was implemented in the TGLF quasi-linear model after improvements to the linear eigenmode accuracy were made. The average fractional error for the 13 scans is 15.6% for ion energy, 14.2% for electron energy, and 35.6% for particle fluxes for TGLF with the SAT2 model and the improved linear eigenmodes.

The spectral shift model for the impact of equilibrium $E \times B$ velocity shear on transport [2] was recalibrated for the SAT2 model. Additional CGYRO scans of the $E \times B$ velocity shear for seven starting cases with different geometry were used for this calibration (30 cases, 23 not in Table 1). It was found that the original model could be simplified to a linear relation between the shift in the radial wavenumber spectrum (eddy tilt) and the $E \times B$ velocity shear. This linear model Eq. 32 fits the CGYRO spectrum of the radial wavenumber shift more accurately than the original non-linear model.

The model for the suppression of electron scale turbulence by zonal flow mixing [3] was also recalibrated to the original multi-scale GYRO simulations [17]. The anisotropy between the radial and poloidal wavenumbers due to geometric effects in the SAT2 model was found to be sufficient to enhance the electron scale fluxes without the need for modification. The coefficient $\alpha_{ZF} = 1.1$ of the zonal flow mixing term (Eq. 35) was hardly changed from unity.

With the recalibration of TGLF, the SAT2 model is now a public option in the TGLF quasi-linear code (SAT_RULE=2). This option automatically includes the geometry modifications and the changes to the electron collision and bounce averaging model.

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Appendix A. Transformation properties of the quasi-linear fluxes

The transformation properties of quasi-linear models have a subtlety that has not previously been discussed in the literature. The gyrokinetic codes typically evolve the Fourier transform of the fields and distribution function. This necessitates the introduction of the interval ΔK_{iy} in the quasi-linear model for the intensity (Eq. 5). It will be shown in this appendix that this factor needs to be defined in a particular way in order for the quasi-linear fluxes to transform in the same way as the non-linear gyrokinetic fluxes for any arbitrary system of normalizations.

For simplicity, consider only the Fourier transform of the angle α defined such that $\mathbf{B} \cdot \nabla \alpha = 0$ [5] and $(\nabla \alpha \cdot \nabla r)_{\theta=0} = 0$:

$$\tilde{\phi}(\alpha) = \sum_{n=-N_{\alpha}}^{N_{\alpha}} \tilde{\phi}_n e^{-in\alpha}$$
(A.1)

$$\tilde{\phi}_n = \int_0^{2\pi} \frac{d\alpha}{2\pi} \tilde{\phi} e^{in\alpha} \tag{A.2}$$

The gyro-kinetic equation does not have an explicit dependence on the mode number "n". The perpendicular wavenumber normalized to the full magnetic field strength in the gyro-average operator (Bessel function) is the physical normalized wavenumber of the turbulence. For a pure ion plasma, with adiabatic electrons, the physical units are the ion temperature, mass and charge, and the magnetic field strength: T_i, m_i, Z_i, B . From these are derived the ion thermal velocity $v_{thi} = \sqrt{2T_i/m_i}$, the ion gyro-frequency, $\Omega_i = Z_i e B/(m_i c)$, and the ion gyro-radius $\rho_i = v_{thi}/\Omega_i$. The normalized poloidal wavenumber that is the argument of the Bessel functions at the outboard midplane for $k_x = 0$ is given by

$$K_{iy} = n\rho_i |\nabla \alpha|_{\theta=0} = \left(\frac{n\rho_i B}{RBp}\right)_{\theta=0}$$
(A.3)

The square amplitude of the Fourier coefficients of the potential fluctuations scales with the interval between ion poloidal wavenumbers (ΔK_{iy}). This property follows from Parseval's theorem

$$\int_{0}^{2\pi} \frac{d\alpha}{2\pi} |\tilde{\phi}|^2 = \sum_{n=-N_{\alpha}}^{N_{\alpha}} |\tilde{\phi}_n|^2 \tag{A.4}$$

A change in the number of mode numbers N_{α} for a fixed range of poloidal wavenumber $K_{iymax} = K_{iy}\Big|_{n=N_{\alpha}}$ requires the Fourier square amplitude to scale like $1/N_{\alpha}$ to maintain the same sum. This scaling is made explicit by defining an intensity function I by the formula:

$$|\tilde{\phi}_n|^2 = \Delta K_{iy} I(K_{iy}) \tag{A.5}$$

In the continuum limit we have

$$\int_{0}^{2\pi} \frac{d\alpha}{2\pi} |\tilde{\phi}|^2 = \sum_{n=-N_{\alpha}}^{N_{\alpha}} \Delta K_{iy} I(K_{iy}) \approx \int_{-K_{iymax}}^{K_{iymax}} dK_{iy} I(K_{iy})$$
(A.6)

The intensity function $I(K_{iy})$ is a function of the ion poloidal wavenumber and so are the linear eigenvalues and quasilinear weights. The ion units are the physical system of units set by the gyro-averaging. Transforming to a new poloidal wavenumber k_{0y} for an arbitrary system of units: T_0, m_0, Z_0, B_0 with $\rho_0 = c\sqrt{m_0T_0}/(Z_0eB_0)$ and

$$k_{0y} = \frac{nq\rho_0 B_0}{rB_{\text{unit}}} \tag{A.7}$$

gives the transformation

$$K_{iy} = \alpha_i \frac{k_{0y}}{|\nabla r|_0} = \beta_i k_{0y} \tag{A.8}$$

where

$$\alpha_i = \sqrt{\frac{2m_i T_i Z_0^2}{m_0 T_0 Z_i^2}} \tag{A.9}$$

In these new units Parseval's Theorem becomes

$$\int_{0}^{2\pi} \frac{d\alpha}{2\pi} |\tilde{\phi}|^{2} = \int_{-k_{0ymax}}^{k_{0ymax}} \beta_{i} dk_{0y} I(\beta_{i} k_{0y})$$
(A.10)

This shows that the correct interval to use in the definition of the quasi-linear intensity Eq. 5 is

$$\Delta K_{iy} = \frac{\alpha_i}{|\nabla r|_0} \Delta k_{0y} \tag{A.11}$$

for any system of units. The CGYRO system of units is given by : $T_0 = T_e, m_0 = m_i, Z_0 = Z_i, B_0 = B_{\text{unit}}$. Thus for the CGYRO units $\alpha_i = \sqrt{2T_i/T_e}$. This factor is

confirmed to be correct by the agreement of the quasilinear fluxes with CGYRO for the scan in T_i/T_e (scan 9).

It can be shown that the product $I_{ky}^{model}WQ_{k_y}$ scales in the expected gyro-Bohm way for the energy flux in changing from one system of units to another. The summation measure Eq. A.11 does not alter the overall scaling of the fluxes since it is the same for any system of units.

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