

## Impact of tungsten charge state bundling on scrape-off layer transport simulations in JET L-mode

H.A. Kumpulainen<sup>1</sup>, M. Groth<sup>1</sup>, M. Fontell<sup>1</sup>, A. Jarvinen<sup>2</sup>, G. Corrigan<sup>3</sup>, D. Harting<sup>3</sup> and JET Contributors\*

<sup>1</sup>*Aalto University, P.O.Box 11000, FI-00076 AALTO, Finland*

<sup>2</sup>*Lawrence Livermore National Laboratory, Livermore, CA 94550, USA*

<sup>3</sup>*CCFE, Culham Science Centre, Abingdon, OX14 3DB, UK*

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**Abstract.** The bundling of the 74 tungsten ion charge states into 6 fluid species in the coupled multi-fluid plasma/kinetic neutral code EDGE2D-EIRENE is shown to decrease the predicted average tungsten charge in the JET main chamber scrape-off layer by up to 40%, compared to mutually consistent predictions using the Monte Carlo impurity transport code DIVIMP and 22 tungsten fluid species in EDGE2D-EIRENE. The tungsten concentrations predicted by DIVIMP in the pedestal region are consistently ~50% higher than by EDGE2D-EIRENE with 6 tungsten fluids under attached divertor conditions, whereas under partially detached conditions the stochastic errors exceed the systematic disagreement between the two codes.

### 1. Introduction

The use of tungsten (W) as the divertor plasma-facing material for ITER, and for the JET ITER-like wall [1], is motivated primarily by its low tritium retention, high melting point, and low erosion [2]. The accumulation of W in the core plasma must be suppressed to ensure that the plasma-cooling effect due to W radiation does not prevent fusion-relevant plasma operation [3]. W transport from the divertor source into the main scrape-off layer (SOL) is critical since it determines the rate at which W enters the core plasma.

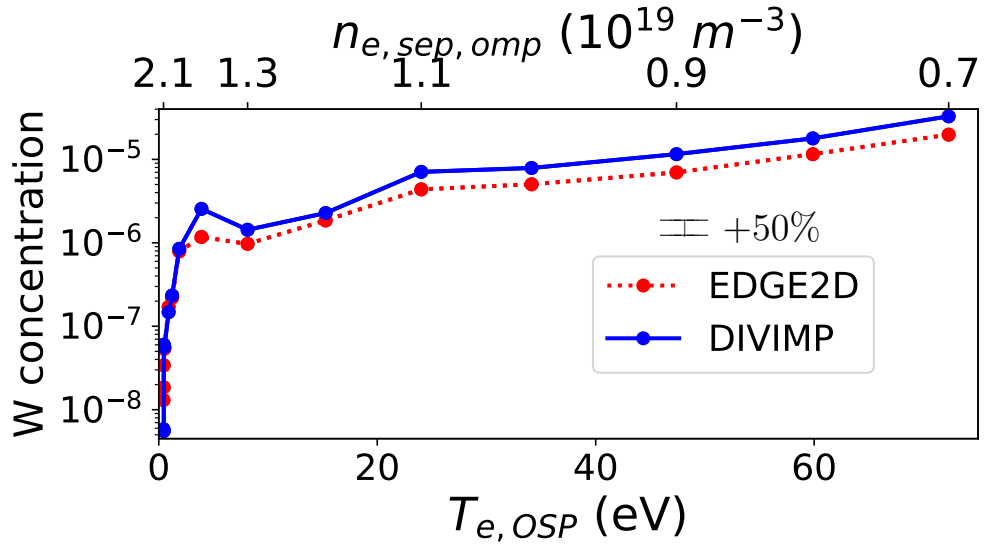
### 2. Simulation setup

The studied plasma scenarios are based on the JET L-mode discharge #81472 (2.5 MA plasma current, 2.5 T toroidal magnetic field) in a vertical HFS/horizontal LFS strike point configuration. The W transport simulations using the coupled 2D multi-fluid plasma/kinetic

neutral code EDGE2D-EIRENE [4, 5] are based on earlier work [6]. The 74 W ion charge states were bundled [7] to 6 fluid species representing charges 1, 2-6, 7-12, 13-22, 23-73 and 74. Particle and energy transport across flux surfaces was treated as purely diffusive with no pinch velocity or cross-field drifts. A range of attached and detached divertor conditions was obtained by altering the prescribed electron density at the LFS mid-plane separatrix while keeping the input power, transport parameters and boundary conditions constant.

In the DIVIMP [8, 9] 2D Monte Carlo simulations, tungsten test particles were injected in a singly ionized state based on the neutral W ionization profile calculated by EDGE2D-EIRENE. The background plasma was imported from EDGE2D-EIRENE. The settings and boundary conditions were selected as close to those used in EDGE2D-EIRENE as possible.

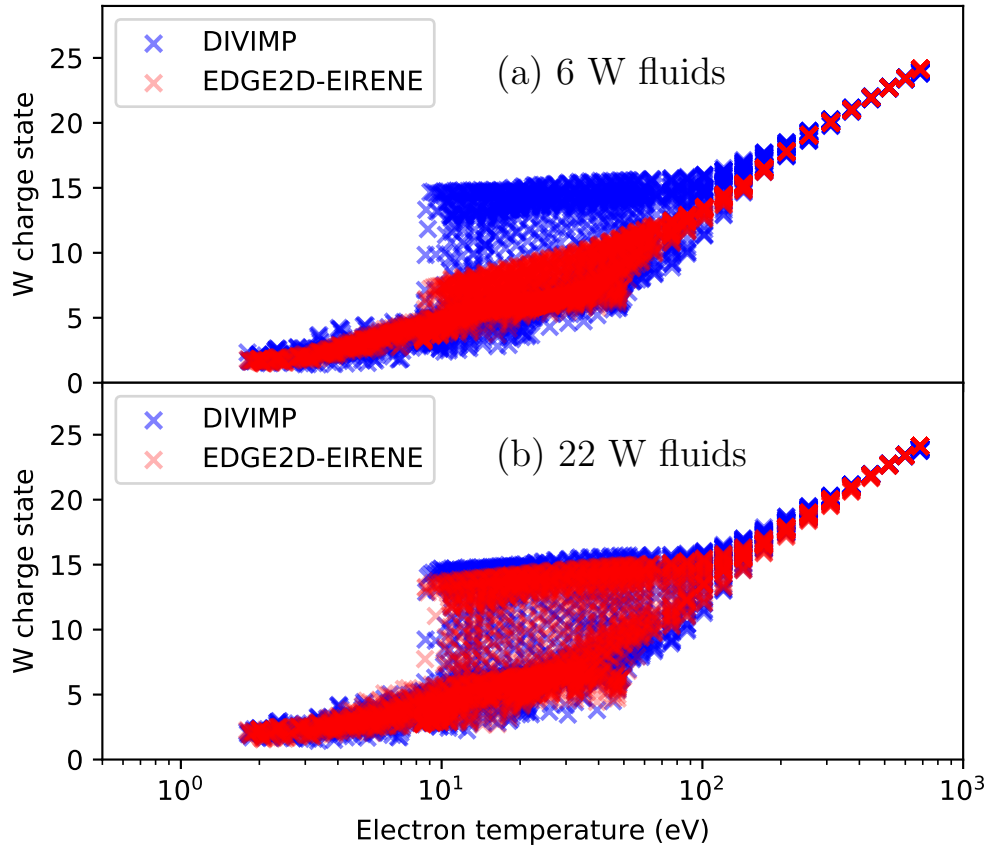
### 3. Comparison of the DIVIMP and EDGE2D-EIRENE predictions



**Figure 1.** Spatial average of the tungsten concentration in the pedestal region (closed flux surfaces at  $r/a > 0.8$ ) predicted by DIVIMP and EDGE2D-EIRENE as a function of electron temperature at the LFS strike point (bottom axis) and electron density at the LFS mid-plane separatrix (top axis).

DIVIMP predicted consistently  $\sim 50\%$  higher W concentration in the pedestal region than EDGE2D-EIRENE under low- and high-recycling divertor conditions (Fig. 1). In partially detached cases with pedestal W concentration below  $10^{-6}$  ( $T_{e,OSP} < 5$  eV), the systematic disagreement was low compared to the Monte Carlo noise in the W source. With very low W concentration ( $< 10^{-7}$ ) in EDGE2D-EIRENE, the error terms in the W particle

and momentum conservation equations became greater than most other terms, essentially rendering the obtained solutions unphysical. Therefore, the conclusions of this work are based on the cases with W concentration  $> 10^{-6}$ .



**Figure 2.** Mean charge state of W ions in each grid cell plotted against electron temperature for DIVIMP and EDGE2D-EIRENE. The 74 W charge states are bundled in EDGE2D-EIRENE into (a) 6 and (b) 22 fluid species.

The bundling of W charge states into 6 fluid species in EDGE2D-EIRENE decreased the average W charge in the upstream SOL region by up to 40% compared to DIVIMP (Fig. 2a). The predicted W charge profiles matched in the divertor ( $1 \text{ eV} < T_e < 60 \text{ eV}$ ,  $1 < Z_W < 6$ ) and in the pedestal region ( $T_e > 100 \text{ eV}$ ,  $Z_W > 13$ ).

Including all the individual W charge states from 1+ up to 20+ as separate fluid species in EDGE2D-EIRENE (the 22-fluid scheme, Fig. 2b), the W charge profiles predicted by EDGE2D-EIRENE and DIVIMP matched within a few percent. The momentum conservation equation, however, was not satisfied due to numerical issues as the EDGE2D-EIRENE runs became unstable. Due to violations of the momentum balance, the 22-fluid scheme did not reproduce the DIVIMP W density profiles despite matching the W charge

profiles.

#### 4. Conclusions

A decrease of 40% in the average charge of W in the main chamber SOL was observed due to the bundling of charge states into 6 fluid species in EDGE2D-EIRENE compared to DIVIMP. This explains the 50% higher W concentration predicted by DIVIMP than by EDGE2D-EIRENE in the pedestal region. Using a more elaborate bundling scheme with 22 W fluids instead of 6, EDGE2D-EIRENE reproduced the W charge profile predicted by DIVIMP within a few percent. However, the increased amount of fluid species in EDGE2D-EIRENE led to larger error terms in the momentum balance. Thus, the 6-fluid W scheme can be considered more appropriate for EDGE2D-EIRENE than the 22-fluid scheme, despite its less realistic W charge profile.

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