Modelling spectral emission from fusion plasmas

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Abstract. The paper adopts initially an historical perspective, commencing from ~ 1970 and has a focus on interests and developments at Culham Laboratory from the time when Nicol Peacock led the UKAEA spectroscopy team. In spite of earlier setbacks, Nicol kept faith in the diagnostic value of quantitative prediction of spectral emissivities in the magnetic confinement fusion domain. So he facilitated and supported continued advance in population modelling in medium density, high temperature, non-LTE, non-stationary plasma on from the original pioneering work of Bates, Kingston and McWhirter (1962) [1]. Calculation of populations of excited states of ions and the distribution between ionisation stages of an element in such plasma requires knowledge of many individual reactions. So precision in spectral prediction is dependent on the availability of such reaction data and carries the additional burden of the generation of such data or its compilation and assessment.

The paper charts a little of the evolution of these models and their data through the seventies and eighties on into this century at Culham, heavily influenced by Nicol. With JET/EP2 and ITER, population structure and ionisation state studies move on, with attention now much focussed on very heavy species, especially tungsten. Modern methods seek to make the atomic physics of such systems manageable for fusion plasma analysis, through techniques such as bundling, superstaging and extended spectral simulation.

Most modern tokamaks have neutral beam heating, so electron collisions are not the end of the population story. Charge transfer between plasma impurity ions and fast neutral hydrogen is the driving reaction for CXS (charge exchange spectroscopy) and impurity ion impact with the fast hydrogen, the driving reaction for BES (beam emission spectroscopy), spectroscopy usually carried out at visible wavelengths. Nicol was interested in all these aspects, although he preferred the VUV. Developments here again recognise ITER and the strengthening of data and models for heavier systems and are included for completeness. But Nicol’s heart was really in spectral line diagnostics, so the paper concludes with the state of efforts to enable easy, universal access to spectral analysis across the scope of Culham activity, of which it is hoped he would approve.

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1 See the Appendix of F. Romanelli et al., Proceedings of the 23rd IAEA Fusion Energy Conference 2010, Daejeon, Korea
INTRODUCTION

It is interesting to look back to the very early days of fusion in the UK from an atomic modelling perspective. Work notes of Thonemann\textsuperscript{2}, an originator of the toroidal confinement approach, include quite detailed considerations of impurity ions in magnetically confined plasmas, such as relaxation of metastable states - which will be returned to later. First efforts to predict electron temperatures in ZETA from line ratios were however misleading because of errors in assumed cross-sections. Pioneering work by Bates, Kingston and McWhirter\textsuperscript{1} on collisional-radiative theory in the early sixties and first R-matrix calculations of cross-sections by Burke had not changed the view of senior management at UKAEA Culham Laboratory in the early seventies that atomic physics and diagnostic spectroscopy was not to be trusted. In any case spectroscopists and atomic physicists at Culham Laboratory were mostly more interested in VUV and soft X-ray observations of the sun’s upper atmosphere from space. Nonetheless Culham had a new divertor tokamak experiment (DITE) coming on-stream which incidentally had molybdenum limiters. So radiant losses from impurity species, including heavy species, mattered and predictive modelling from atomic physics was required and tolerated. It was then, in the early seventies, that one of us (Summers) first became involved in magnetic confinement fusion at Culham. In fact this was with Nicol Peacock, who had kept faith with atomic modelling for plasmas and wished for some calculations on ionisation state and radiated power of species in DITE. So, from that point in time, a connection was established with Nicol which has continued on through the eighties into this century with the JET Joint Undertaking. It was of course clear that lack of accurate knowledge of electron-ion reaction rates limited our work and fortunately, in the late eighties, another of us (Badnell) started a crucial, long-term focus on precision cross-sections which continues to this day. To complete the story, another of us (O’Mullane) a research student with Nicol Peacock in the nineties continues our efforts into ITER and the future and in fact wrote Nicol’s last paper with him\textsuperscript{2}. It must be said that Nicol did not stand still, but was always enthused by the latest ideas. So the rest of this paper is not a historical work, but rather a tribute, showing where the engagement with Nicol has led us to now.

POPULATION STRUCTURE AND IONISATION STATE

In our Cambridge ivory tower in the early seventies, Nicol appeared one day excited about titanium ions and wanting to model the low level population structure so that he could predict the emissivities of the associated VUV lines. This was an interesting departure, since at that time our main concern was dielectronic recombination and its sensitivity to finite plasma density - a very highly excited Rydberg level problem. Also, Nicol was concerned about the emission of OVII lines in the recombining environment of the peripheral DITE plasma which seemed to reveal differential behaviour of ground and metastables in radial transport. Again this was intriguing, since early errors were made in effective dielectronic recombination from neglect of abundant metastable re-

\textsuperscript{2} Private communication and notebook archives from R. W. P. McWhirter
combining target ions. Later, at the JET Joint Undertaking in the middle eighties, related questions, associated with impurity influx from localised surfaces contacted by the plasma, arose. Spectral observations of such inflowing ions are conveniently performed in the visible and so arise from excited levels in higher n-shells for light element ions. A typical transition is OII (2s²2p²3p⁴P − 2s²2p²3s⁴S) whose emissivity is driven from the 2s²2p²3s⁴S ground primarily. But the 2s²2p²3p⁴P level is relatively long-lived, since it cannot decay to the ground. It is therefore susceptible to collisional re-distribution at tokamak densities and also to augmentation by cascade from higher n-shells. Although JET was a machine following a light element first wall strategy (beryllium and carbon), influx of more complex heavier elements such as Cr (JET antennae screens) and Mo (DITE and ALCATOR) were of interest. These motivations set in train the pattern of excited population and ionisation state modelling which, in a more refined form, we use and continue to develop today. This can be summarised as trying to address the whole excited population structure of an ion, ground to continuum, recognizing metastability, including all relevant collisional processes, adopting collisional-radiative principles on timescale ranking to obtain ‘effective coefficients’, and using varying resolutions (in a shell, term or level sense) and population grouping to make the problem computationally manageable (see figure 1). The realisation of these principles in codes, together with all the data, fundamental and derived, necessary to make it practical and relevant became
ADAS, the Atomic Data and Analysis Structure\textsuperscript{3}. The ADAS Project is grateful to Nicol for his encouragement in this area over many years.

To elaborate a little on the present state of affairs, ADAS population modelling, at its highest precision, has been targeted at the ions of the elements hydrogen to neon. The details of the methods are described in Summers \textit{et al.} (2006) \cite{3}. A simplified schematic is shown in figure 1 (but see also Summers \textit{et al.} (2011) \cite{4}). Comprehensive derived effective coefficient data are available to support modelling of the separate dominant metastables of light elements in plasma transport equations and prediction of associated spectral emission. Periodic updates of the data occur in response to improved calculations of electron-impact cross-sections between low levels (see figure 1).

\textbf{FIGURE 2.} Emission lines observed by SUMER, CDS and EIS in the joint solar observations JOP220/HOP109 on 17\textsuperscript{th} April 2009 \cite{5} together with the instrument wavelength coverages. The temperatures of line formation have been shown, in order to indicate the solar atmosphere layer from which the indicated lines arise. The spectral ranges of JET spectrometers becoming operational in the EP2 extended performance upgrade and relevant divertor zones are also indicated for comparison.

However, new differential emission measure (DEM) analyses in the lower temperature solar chromosphere/transition region (Giunta (2011) \cite{5}), and of course preparations for ITER in fusion, have compounded pressure to extend the range of species to argon and possibly iron. This is further justified by new state-resolved dielectronic recombination, ionisation cross-section and R-matrix excitation cross-section calculations which provide underpinning fundamental data. Silicon is now complete and magnesium will soon follow. Analyses using these data do indicate that all of the ingredients summarised in the previous paragraph matter. The studies confirm that truncation of the population structure at a set of low levels and failure to include the finite-density suppression of

\textsuperscript{3} URL: http://www.adas.ac.uk
dielectronic recombination, even with good low-level fundamental data, does lead to mis-interpretation [5]. In such development, fusion and astrophysics can be mutually supportive. Figure 2 shows our current focus of interest in solar observations and the relationship to divertor spectrometry and species at Culham laboratory (JET and MAST). The derived atomic data required for DEM analysis are contribution functions as illustrated in figure 3, but their ingredients (effective emissivity, ionisation and recombination coefficients) allow equivalent prediction in non-stationary transport regimes and transients of both the solar atmosphere and tokamak divertors (except that the tokamak evolution is about one thousand times faster).

It is though the change of emphasis in population and emissivity modelling from light to heavy elements which has been the main pre-occupation. The methods being used for heavy element ions were outlined in Summers et al. (2007) [6] and in O’Mullane et al. (2009) [7] and since then they have become more refined, systematised and tested. Superstaging has now been carried into 2-D transport modelling enabling studies of tungsten (Strachan et al. (2010)[8]). This has encouraged efforts to improve the baseline approximations used in ADAS and to target selected ions at the highest precision. Figure 4 shows a partition which isolates the ions W$^{+20}$ and W$^{+44}$ from the grosser superstages.

Figure 5 shows the level counts used for representation of the power, drawing attention to the above specific stages. Within the JET scope of core electron temperature, W$^{+44}$ stands out as a somewhat simpler system with predicted spectral emission in the range of the KX1 soft X-ray spectrometer. First relativistic, R-Matrix, radiation damped cross-section calculations have been carried out by Ballance and Griffin (2007) [9], but omit the observable lines which are 3d – 4f core excited. Further very large calculations are in preparation to explore this. The W$^{+20}$ ion with a partially filled f-shell is a formidable system with a quasi-continuum of emission, yet marks (to lower ionisation stages) the divertor relevant stages. In ionisation, the CADW approximation of Loch et al. [10] (see also Loch et al. [11]) raises our baseline, but measurements by Borovik et
FIGURE 4. Tungsten equilibrium ionisation balance in superstage partitioning with isolation of $W^{+20}$ and $W^{+44}$ for more detailed spectral studies.

FIGURE 5. Transition counts for tungsten ionisation stages in baseline modelling. Systems such as $W^{+44}$ have fewer, more intense, lines - some in the JET soft X-ray spectrometer range. $W^{+20}$ by contrast, with a half-open f-shell, presents difficulties for computation and validation of atomic data.

al. (2011) [12] still indicate discrepancies even at quite high charge states from REDA-like structures. AUTOSTRUCTURE raises our baseline for dielectronic recombination (see Badnell et al. (2003) [13]), and is being pushed towards more complex systems (Badnell et al. (2011) [14]). These comparative calculations suggest a varying bundling
and resolution approach along with approximations such as BBGP [13] is the way forward. Such an approach can probably make manageable both the low temperature and high temperature needs. Again the measurements (Schippers et al. (2011) [15]) show that very large low energy resonances can influence the high temperature dielectronic rates. Our current work moves forward in all these aspects, but in a strongly targeted manner driven by observability as indicated above.

**NEUTRAL BEAMS AND CHARGE EXCHANGE SPECTROSCOPY**

Nicol was not too enthused by charge exchange spectroscopy (CXS) or at least by that conducted at visible wavelengths and directed at inferring concentrations. Partly this reflected his VUV and shorter wavelength preference, but also because the observations are of emission from sub-dominant levels of the receiver with rather a long trail of atomic modelling involving, sometimes, poorly known cross-sections. He would therefore be shocked by figure 6, an estimate designed to examine how much tungsten active signals could affect fast helium concentration deduction from charge exchange spectroscopy.

However CXS certainly was and remains a subject of interest and much discussion. State selective charge exchange data from neutral hydrogen ground state and excited donors to fully stripped light impurity nuclei as receivers has improved. Recent studies consolidate and extend the coverage with precision methods such as CCMO for B\(^{+5}\) (Guzman et al. (2010)[16]) and CCAO for Be\(^{++}\)(Igenbergs et al. (2009) [17]) and N\(^{+7}\) (Igenbergs (2011) - private communication). Extension to argon is the prize because of ITER relevance and also because it allows (by simultaneous observation of Ar\(^{+15}\), Ar\(^{+16}\) and Ar\(^{+17}\) CXS emission in the visible) a direct connection with impurity transport. Differences and discontinuities between methods (CTMC, modified CTMC and CCMO) in the key ~ 50keV/amu region are an anxiety. Calculations by Igenbergs (private communication) in CCAO approximation will hopefully clarify the situation. Meanwhile, the existing database does allow the preparation of universal semi-empirical z-scaled
parametric forms for the state selective data (Foster (2008) [18]) which have been included in ADAS and used for extrapolation to high Z (see figure 6). Further studies on CXS of a Kr$^{+36}$ receiver in the modified CTMC approach (Illescas (2011) - private communication) and the Ar$^{+18}$ revision will strengthen the universal parametric form. The conversion of such data to spectral line emissivities uses the same form of models as discussed in figure 1.

SPECIAL FEATURES AND SPECTRAL ANALYSIS

**FIGURE 7.** ADAS605 processing screen: allows interactive manipulation of the chosen feature via custom control widgets in the right hand panel, with graphical output in the left panel.

The now extensive population and ionisation state models of ADAS, which span most fusion plasma environments and species to some measure, combined with their delivery of derived coefficients in precisely defined data formats, provide a basis for spectral diagnostics. Helium/lithium-like satellite lines, Balmer decrements, Zeeman pattern lines and so on are just groupings of related lines exploited as special features because of their special sensitivity. Ad hoc models and codes abound to expedite such exploitation. The comprehensive character of ADAS has encouraged us to explore and expand the atomic physics link so that the special features may be realised as a universal mathematical/computational construct for use in spectral modelling and fitting. In close association with ADAS, one of us (Nicholas) [19] has sought to empower special feature analysis by implementing such generalised computational structures — AFG (ADAS Feature Generator) and FFS (Framework for Feature Synthesis). These allow both a pedagogical insight into the capabilities of each special feature as well as practical execution of optimised spectral fitting and plasma parameter extraction. The methods, based on object-oriented programming, are universal including aspects such as self-generating graphical user interfaces and an algebra of parametric feature
creation. Special features were close to Nicol’s heart and so this paper concludes with some illustrations of our latest methods on selected JET spectra.

**FIGURE 8.** C III ($1s^22s3s^3S - 1s^22s3p^3P$) multiplet at $\sim 465$ nm as recorded by KS8 at JET, pulse #75898. The in-board and out-board emission, differentially perturbed by the magnetic field and flow, along a mid-plane, horizontal line-of-sight, form the observational feature. The fitted FFS model is shown in red, with the low-field and high-field side Zeeman component features indicated in green and blue, respectively. The fit residual is shown in the lower plot.

The main drive behind AFG is to ease access to the ADAS special feature routines such that they are easily incorporated into any external modelling code (such as FFS). It is possible to do this through a series of simple commands — now common to all of the ADAS special features. However, it was thought that AFG could be made even more accessible via a graphical user interface (GUI); this code is known as ADAS605. It has been designed to use the AFG API intensively, such that the interface is highly dynamic. Its appearance is very much dependent upon the feature under consideration. As illustration of a Zeeman feature is shown in figure 7.

Theoretical representations of experimental spectra can be considered to be constructs from a set of model elements; from simpler mathematical line shapes to more complex special feature representations from ADAS. One can also define a set of ‘operator elements’ which can apply a function to the results of other elements, in some cases the function will bring together combinations of element results to represent the spectra in its entirety. Programmatically, each of the elements can be represented by an object that provides a method for performing calculation of that component of the spectra. This is the approach taken by the Framework for Feature synthesis (FFS). In order to specify the construct for an arbitrarily complex model spectra, a LISP-like model definition language (MDL) has been set up. Details of initial user model optimising, allowing substitution of analytic derivatives and hence very substantial fitting speed-up
are beyond this short paper (see Nicholas (2011)[19]). An illustration from JET is shown in figure 8.

CONCLUSIONS

Nicol was a fusion and spectroscopy enthusiast who freely shared information, ideas and results. He facilitated engagement between experimental spectroscopists, theoretical atomic physicists and plasma physicists. Also, he had an outward looking perspective from Culham Laboratory which allowed external university scientists to become fully engaged. We believe this was very fruitful as indicated by this brief survey of what was started and where it has got to now. His legacy at Culham is evident. ITER is our next target in which Nicol would surely have concurred.

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