Plasma Chemistry and Kinetics in Low Pressure Oxygen Discharges: The significance of metastable states

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Outline

- A. Global (volume averaged) chemistry models
 - A.1 The oxygen discharge
 - chlorine dilution
- B. 1D particle-in-cell/Monte Carlo collision simulation
 - B.1 Capacitively Coupled Oxygen Discharge at 13.56 MHz
 Voltage Source pressure dependence
- C. Summary





A. Global (volume averaged) chemistry models





Global (volume averaged) chemistry models

- The main idea of a global model is to generate a model that encompasses a large number of reactions in order to model a processing plasma with a limited computing power by neglecting the complexity which arises when spatial variations are considered
- Thus the model does not describe spatial distribution but captures scalings of plasma parameters with control parameters
- The model allows us to investigate various phenomena, such as the effects of excited species, negative ions and particular reactions on the overall discharge





A.1 Global (volume averaged) chemistry models – oxygen





The oxygen discharge

- The oxygen discharge is of significance in various materials processing applications including
 - etching of polymer films
 - ashing of photoresist
 - oxidation
- The oxygen chemistry is complicated due to the presence of metastable atomic and molecular species
- It is in particular the two low lying metastable molecular states designated by $a^1 \Delta_g$ and $b^1 \Sigma_g^+$, which are located 0.98 and 1.627 eV above the ground state, respectively
- It is well established that collisions with these metastable states have in many cases larger cross sections and thus higher reaction rates than corresponding collisions with the ground state molecule

- A steady state global (volume averaged) model was developed for the oxygen discharge
- The following species are included
 - electrons
 - the ground state oxygen molecule $O_2(X^3\Sigma_g^-, v = 0)$,
 - The metastable oxygen molecules $O_2(a^1\Delta_g)$, $O_2(b^1\Sigma_g^+)$ and the metastable Herzberg states $O_2(A^3\Sigma_u^+, A'^3\Delta_u, c^1\Sigma_u^-)$
 - the ground state oxygen atom O(³P)
 - the metastable oxygen atom O(¹D)
 - the negative oxygen ions O⁻ and O⁻₂
 - the positive oxygen ions O⁺ and O⁺₂
 - Ozone O₃ and its ions O₃⁺ and O₃⁻
- The content of the chamber is assumed to be nearly spatially uniform and the power is deposited uniformly into the plasma bulk

■ The particle balance equation for a species *X* is given

$$\frac{\mathrm{d}n^{(X)}}{\mathrm{d}t} = 0 = \sum_{i} R_{\mathrm{Generation},i}^{(X)} - \sum_{i} R_{\mathrm{Loss},i}^{(X)}$$

where $R_{\text{Generation},i}^{(X)}$ and $R_{\text{Loss},i}^{(X)}$, respectively, are the reaction rates of the various generation and loss processes of the species X

■ The power balance equation, which equates the absorbed power P_{abs} to power losses due to elastic and inelastic collisions and losses due to charged particle flow to the walls is given as

$$\frac{1}{V} \left[P_{\text{abs}} - eVn_{\text{e}} \sum_{\alpha} n^{(\alpha)} E_{\text{c}}^{(\alpha)} k_{\text{iz}}^{(\alpha)} - eu_{\text{B0}} n_{\text{i}} A_{\text{eff}} (E_{\text{i}} + E_{\text{e}}) \right] = 0$$

For the edge-to-center positive ion density ratio we use

$$h_{\ell} \simeq \left[\left(\frac{0.86}{(3 + \eta L/2\lambda_{i})^{1/2}} \frac{1}{1 + \alpha_{0}} \right)^{2} + h_{c}^{2} \right]^{1/2}$$

$$h_{R} \simeq \left[\left(\frac{0.8}{(4 + \eta R/\lambda_{i})^{1/2}} \frac{1}{1 + \alpha_{0}} \right)^{2} + h_{c}^{2} \right]^{1/2}$$

where $\alpha_0 \approx (3/2)\alpha$ is the central electronegativity, $\eta = 2T_+/(T_+ + T_-)$ and

$$h_{\rm c} \simeq \left[\gamma_-^{1/2} + \gamma_+^{1/2} \left[n_*^{1/2} n_+ / n_-^{3/2} \right] \right]^{-1} \quad {\rm and} \quad n_* = \frac{15}{56} \frac{\eta^2}{k_{\rm rec} \lambda_{\rm i}} v_{\rm i}$$

is based on a one-region flat topped electronegative profile

$$\gamma_- = T_e/T_-$$
 and $\gamma_+ = T_e/T_+$



- The electron energy distribution function (EEDF) is usually assumed to be Maxwellian
- We can also assume a general electron energy distribution

$$g_{\rm e}(E) = c_1 E^{1/2} \exp\left(-c_2 E^x\right)$$

$$c_1 = \frac{1}{\langle E \rangle^{3/2}} \frac{\left[\Gamma(\xi_2)^{3/2}\right]}{\left[\Gamma(\xi_1)^{5/2}\right]} \quad \text{and} \quad c_2 = \frac{1}{\langle E \rangle^x} \frac{\left[\Gamma(\xi_2)\right]}{\left[\Gamma(\xi_1)\right]^x}$$

where
$$\xi_1 = 3/2x$$
 and $\xi_2 = 5/2x$

■ Here x = 1 and x = 2 correspond to Maxwellian and Druyvesteyn electron energy distributions, repectively

$$g_{\mathrm{p}}(E) = \frac{g_{\mathrm{e}}(E)}{E^{1/2}}$$





■ The diffusional losses of the neutral oxygen atoms (ground state and metastable) to the reactor walls are given by

$$k_{\text{O,wall}} = \left[\frac{\Lambda_{\text{O}}^2}{D_{\text{O}}} + \frac{2V(2 - \gamma_{\text{rec}})}{Av_{\text{O}}\gamma_{\text{rec}}}\right]^{-1} \text{ s}^{-1}$$

- D_O is the diffusion coefficient for oxygen atoms
- $v_{\rm O} = (8eT_{\rm g}/\pi m_{\rm O})^{1/2}$ is the mean O velocity
- $\ \ \, \ \, \ \, \gamma_{\rm rec}$ is the wall recombination coefficient for neutral oxygen atoms on the wall surface
- lacktriangle Λ_O is the effective diffusion length of neutral oxygen atoms

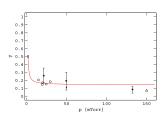
$$\Lambda_{\rm O} = \left[\left(\frac{\pi}{L} \right)^2 + \left(\frac{2.405}{R} \right)^2 \right]^{-1/2}$$

■ The wall recombination probability, γ_{rec} , is a very important quantity in all low pressure molecular discharges



Surface recombination

- The pressure dependence on the wall recombination coefficient was achieved by fitting all the available data for stainless steel surfaces
- For anodized aluminium reactor walls the recombination coefficient is assumed to be a constant γ_{rec} = 0.06 (Guha et al (2008a))
- The same wall recombination coefficient was used for O(¹D) as no data is available



The wall recombination coefficient for oxygen atoms on stainless steel surfaces depends on pressure through

 $\gamma_{\text{rec}} = 0.1438 \exp(2.5069/p)$ p > 2 mTorr

$$\gamma_{\text{rec}} = -0.25p + 1$$
 $p < 2 \text{ mTorr}$



A.1.1 Particle densities

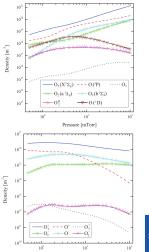




Particle densities

- The dominant species is the oxygen molecule in the ground state O₂(X³Σ_g) followed by the oxygen atom in the ground state O(³P)
- The singlet metastable states $O_2(a^1\Delta_g)$ and $O_2(b^1\Sigma_g^+)$ and the metastable atom $O(^1D)$ are also present in the plasma in significant amounts
- a cylindrical stainless steel chamber
 radius R = 15 cm and length L = 30 cm P_{abs} = 500 W

Toneli et al., J. Phys. D, 48 (2015) 325202

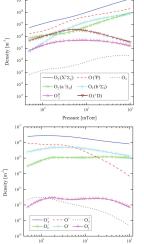


Pressure [mTorr]



Particle densities

- The $O_2(b^1\Sigma_g^+)$ density overcomes the $O_2(a^1\Delta_g)$ density in the pressure range from 2.5 to 80 mTorr
- The O₂⁺ ions are in majority among the positive ions
- The O⁺ density has a sharp decrease for pressures above 4 mTorr
- The ratio [O⁻]/[O₂⁻] is 5.3 at 1 mTorr, and 1.3 at 100 mTorr
- a cylindrical stainless steel chamber
 radius R = 15 cm and length L = 30 cm P_{abs} = 500 W



Pressure [mTorr]



Creation of metastable $O_2(a^1\Delta_g)$ molecules

■ Electron impact excitation (17)

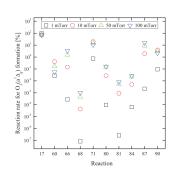
$$\mathrm{e} + \mathrm{O}_2(\mathrm{X}^3\Sigma_g^-) \longrightarrow \mathrm{O}_2(\mathrm{a}^1\Delta_g) + \mathrm{e}$$

has 100 % contribution at 1 mTorr and 68.2 % at 100 mTorr

■ The role of quenching of the metastable atom O(¹D) to create the metastable oxygen molecule (71)

$$O(^{1}D)+O_{2}(X^{3}\Sigma_{g}^{-})\longrightarrow O_{2}(a^{1}\Delta_{g})+O(^{3}P)$$

reaches a maximum value, 20.1 % contribution at 12.5 mTorr, and 10.5 % at 100 mTorr



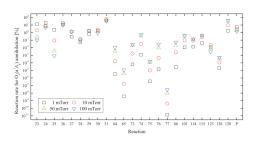
Toneli et al., J. Phys. D, 48 (2015) 325202







Destruction of metastable $O_2(a^1\Delta_g)$ molecules



■ Electron impact dissociation (31,24)

$$\begin{split} &e + \mathrm{O}_2(\mathrm{a}^1 \Delta_g) \longrightarrow \mathrm{O}(^1\mathrm{D}) + \mathrm{O}(^3\mathrm{P}) + \mathrm{e} \\ &e + \mathrm{O}_2(\mathrm{a}^1 \Delta_g) \longrightarrow \mathrm{O}(^3\mathrm{P}) + \mathrm{O}(^3\mathrm{P}) + \mathrm{e} \end{split}$$

are the most important channels for destruction of the metastable $O_2(a^1\Delta_g)$ molecules, 57.2 % contribution at 1 mTorr and 38.1 % at 100 mTorr



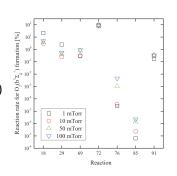
Creation of metastable $O_2(b^1\Sigma_g)$ molecules

■ The most important contributor to the formation of the metastable oxygen molecule $O_2(b^1\Sigma_g)$ is reaction 72

$$O(^{1}D)+O_{2}(X^{3}\Sigma_{g}^{-})\longrightarrow O_{2}(b^{1}\Sigma_{g})+O(^{3}P)$$

which has 75.7 % contribution at 1 mTorr and 93.2 % at 100 mTorr

■ This reaction has been suggested in the past to be a major contributor to the formation of $b^1\Sigma_g^+$ state in the atmosphere



Toneli et al., J. Phys. D, 48 (2015) 325202







Destruction of metastable $O_2(b^1\Sigma_g)$ molecules

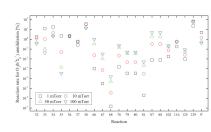
- Wall quenching is an important loss process for $O_2(b^1\Sigma_q)$ (129)
- Electron impact dissociation (38,34)

$$e + O_2(b^1\Sigma_g) \longrightarrow O(^1D) + O(^3P) + e$$

$$e + O_2(b^1\Sigma_g) \longrightarrow O(^3P) + O(^3P) + e$$

are the most important channels for destruction along with electron impact ionization (33)

$$e + O_2(b^1\Sigma_g) \longrightarrow O_2^+ + 2e$$



Toneli et al., J. Phys. D, 48 (2015) 325202







Destruction of the negative ion O-

 At low pressure that the electron impact detachment (54),

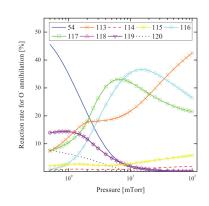
$$e + O^- \longrightarrow O(^3P) + 2e$$

is the main contributor

■ Ion-ion mutual neutralization is also important at low pressure, note that for O⁻ + O⁺₂ we use

$$k = 2.6 \times 10^{-14} (300/T_{\rm g})^{0.44} \, {\rm m}^3/{\rm s}$$

 As pressure increases, the contributions of these reactions become negligible





Destruction of the negative ion O-

 As the pressure increases charge exchange (113)

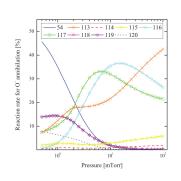
$$\mathrm{O}^- + \mathrm{O}_2(\mathrm{X}^3\Sigma_g^-) \longrightarrow \mathrm{O}(^3\mathrm{P}) + \mathrm{O}_2^-$$

and detachment by collision with O(³P) (117)

$$\mathrm{O}^- + \mathrm{O}(^3\mathrm{P}) \longrightarrow \mathrm{O}_2(\mathrm{X}^3\Sigma_g^-) + \mathrm{e}$$

have increased contribution as well as detachment by collision with $O_2(b^1\Sigma_g)$ (reaction 116)

$$O^-+O_2(b^1\Sigma_g) \longrightarrow O(^3P)+O_2(X^3\Sigma_g^-)+e$$



Toneli et al., J. Phys. D, 48 (2015) 325202







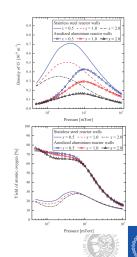
A.1.2 Influence of chamber wall and EEDF





Influence of chamber wall and EEDF

- The parameter *x* defines the shape of the electron energy distribution
 - $\mathbf{x} = 0.5$ is concave or bi-Maxwellian
 - $\mathbf{x} = 1$ is Maxwellian distribution
 - $\mathbf{x} = 2$ is Druyvesteyn distribution
- For anodized aluminium reactor walls, the recombination coefficient for oxygen atoms at the walls is assumed to be a constant $\gamma_{\rm O}$ = 0.06 based on the measurements of Guha et al. (2008a)
- The chamber wall has a significant influence on the dissociation fraction



A.1.3 Chlorine dilution



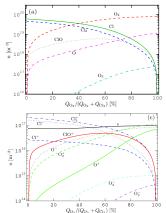


Chlorine dilution - Particle densities

- The chlorine-oxide molecule CIO and its ion CIO⁺ peak when CI₂ and O₂ flowrates are roughly equal
- The $O_2(a^1\Delta_g)$ density is about 9 10 % of the total O_2 density
- The electron density increases about 30 % between pure chlorine and pure oxygen discharge

Thorsteinsson and Gudmundsson, Plasma Sources Sci.

Technol., 19 055008 (2010)

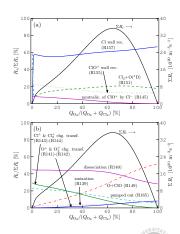


A cylindrical stainless steel chamber L = 10 cm and R = 10 cm p = 10 mTorr and $P_{abs} = 500$ W



Oxygen dilution - Particle densities

- The total rate for creation and loss of CIO molecules is at maximum when the oxygen content is 65%.
- Wall recombination of CI molecules, is the dominating pathway for creation of CIO molecules
- The bulk processes and recombination of CIO⁺ ions at the wall account for roughly 33–43% of the total rate for CIO creation, combined



Thorsteinsson and Gudmundsson, Plasma Sources S

Technol., 19 055008 (2010)_





A.1.4 Summary





Summary

- A global model of O_2 , and Cl_2/O_2 discharges has been applied to explore the creation and destruction of the singlet metastable molecules $O_2(a^1\Delta_g)$, $O_2(b^1\Sigma_g^+)$
- The singlet delta state $O_2(a^1\Delta_g)$ is created mainly by lectron impact excitation

$$e + O_2(X^3\Sigma_g^-) \longrightarrow O_2(a^1\Delta_g) + e$$

■ The singlet delta state $O_2(b^1\Sigma_g^+)$ is created mainly by

$$O(^{1}D) + O_{2}(X^{3}\Sigma_{g}^{-}) \longrightarrow O_{2}(b^{1}\Sigma_{g}) + O(^{3}P)$$

At low pressure that the electron impact detachment

$$e + O^- \longrightarrow O(^3P) + 2e$$

is the main contributor to the loss of O-

■ The CIO molecule is mainly created by recombination at the discharge wall





B. 1D particle-in-cell/Monte Carlo collision simulation





Outline

- The 1D particle-in-cell/Monte Carlo collision simulation
 - The oxygen discharge
 - Capacitively Coupled Oxygen Discharge at 13.56 MHz Voltage Source – pressure dependence
 - Summary





The 1D particle-in-cell/Monte Carlo collision simulation





The oopd1 1d-3v PIC/MCC code

- In particle-in-cell simulation the plasma is represented as a collection of macroparticles
- Each macroparticle is a charged "cloud" representing many real charged particles
- Each macroparticle has the same charge-to-mass ratio (q/m) as the real charged particle
- Equations of motion are solved for each macroparticle
- The electric and magnetic fields are calculated self-consistently using charge densities and currents produced by the macroparticles





The oopd1 1d-3v PIC/MCC code

- We use the oopd1 (objective oriented plasma device for one dimension) code to simulate the discharge
- The oopd1 code was originally developed at the Plasma Theory and Simulation Group at UC Berkeley
- It has 1 dimension in space and 3 velocity components for particles (1d-3v)
- The oopd1 code is supposed to replace the widely used xpdx1 series (xpdp1, xpdc1 and xpds1)
- It is developed to simulate various types of plasmas, including processing discharges, accelerators and beams
 - Modular structure
 - Includes relativistic kinematics
 - Particles can have different weights





B.1. The oxygen discharge





The oxygen discharge

- We consider a discharge that consists of:
 - electrons
 - the ground state oxygen molecule $O_2(X^3\Sigma_g^-)$
 - the metastable oxygen molecule $O_2(a^1\Delta_g)$
 - the metastable oxygen molecule $O_2(b^1\Sigma_g)$
 - the ground state oxygen atom O(³P)
 - the metastable oxygen atom O(¹D)
 - the negative oxygen ion O⁻
 - the positive oxygen ions O⁺ and O⁺₂
- We apply a global model¹ beforehand to calculate the partial pressure of the various neutrals

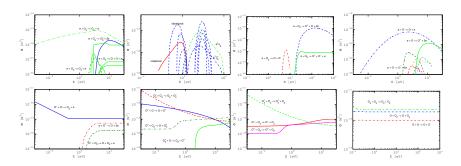






Thorsteinsson and Gudmundsson, Plasma Sources Sci. Technol., 19 055008 (2010)

The oxygen discharge



■ The reaction set for the oxygen is comprehensive and for this study includes 67 reactions

Gudmundsson et al., Plasma Sources Sci. Technol., 22 035011 (2013), 24 035016 (2015) and 25 055002 (2016)





B. 1. 1. Capacitively Coupled Oxygen Discharge at 13.56 MHz – pressure dependence – including $O_2(a^1\Delta_a)$





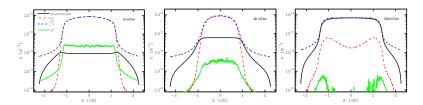
We apply a voltage source with a single frequency

$$V(t) = V_{\rm rf} \sin(2\pi f t)$$

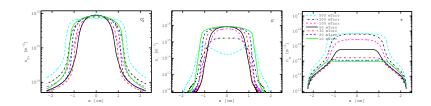
- The electrodes are circular with a diameter of 14.36 cm
- The gap between the electrodes is 4.5 cm
- We set V_{rf} = 222 V and f = 13.56 MHz
- The neutrals (O_2 and O) are treated as background gas at $T_g = 300$ K with a Maxwellian distribution
- The dissociation fraction and the metastable fraction is found using a global model
- The pressure is varied from 10 500 mTorr







- For a parallel plate capacitively coupled oxygen discharge at 50 mTorr with with a gap separation of 4.5 cm by a 222 V voltage source at 13.56 MHz
 - O₂⁺-ion density profile
 - O⁺-ion density profile
 - O⁻-ion density profile
 - electron density profile

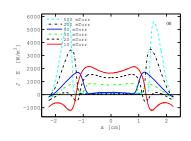


- The sheath width decreases as the pressure is decreased in the pressure range from 50 mTorr to 10 mTorr
- The sheath widths are largest at 50 mTorr
- As the pressure is increased from 50 mTorr up to 500 mTorr the sheath width decreases
- This agrees with what has been observed experimentally in the pressure range 40 375 mTorr

Mutsukura et al. (1990) JAP 68 2657 and van Roosmalen et al. (1985) JAP 58 653



- The electron heating profile $J_e \cdot E$
- In the pressure range 50 500 mTorr the electron heating occurs almost solely in the sheath region
- As the pressure is decreased the Ohmic heating contribution in the plasma bulk increases and sheath heating decreases

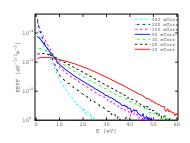


Gudmundsson and Ventéjou (2015) JAP 118 153302





- At low pressure the EEPF is convex, the population of low energy electrons is relatively low
- As the pressure is increased the number of low energy electrons increases and the number of higher energy electrons (> 10 eV) decreases
- Thus the EEPF develops a concave shape or becomes bi-Maxwellian as the pressure is increased



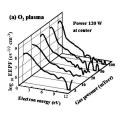
Gudmundsson and Ventéjou (2015) JAP 118 153302



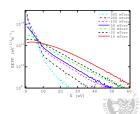




- Our results agree with the measurements of Lee et al.
 (2010) which explored experimentally the evolution of the EEPF with pressure in a capacitively coupled oxygen discharge in the pressure range 3 100 mTorr
- They find that the EEPF became more distinctly bi-Maxwellian and the density of low energy electrons increases as the gas pressure is increased



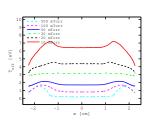
Lee et al. (2010) PRE 81 046402



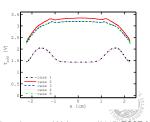


- The effective electron temperature drops as the pressure is increased
- When the metastable singlet oxygen molecule $O_2(a^1\Delta_g)$ is added to the discharge model the effective electron temperature drops, in particular in the electronegative core due to detachment by the metastable $O_2(a^1\Delta_g)$ molecule

$$O^- + O_2(a^1 \Delta_a) \rightarrow products$$

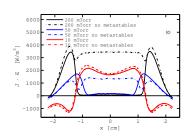


Gudmundsson and Ventéjou (2015) JAP 118 153302



Gudmundsson and Lieberman (2015) PSST 24 0350

- At 10 mTorr excluding the metastable states in the simulation has very small influence on the heating mechanism
- At 50 mTorr the metastable states have a significant influence on the heating mechanism
- The role of the metastables is even more significant at 200 mTorr



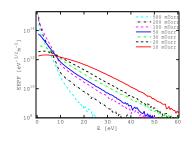
Gudmundsson and Ventéjou (2015) JAP **118** 153302 Gudmundsson and Lieberman (2015) PSST **24** 035016







- At low pressure the EEPF is convex and develops a concave shape or becomes bi-Maxwellian as the pressure is increased
- These results contradict what is commonly found for the capacitively coupled argon discharge where the EEPF evolves from being concave at low pressure to being convex at high pressure



Gudmundsson and Ventéjou (2015) JAP 118 153302





B. 1. 2. Capacitively Coupled Oxygen Discharge at 13.56 MHz – pressure dependence – including $O_2(\mathbf{a}^1\Delta_g)$, $O_2(\mathbf{b}^1\Sigma_g)$ and $\gamma_{\rm see}(E)$



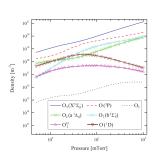


■ It has been known for decades that the metastable oxygen molecule $O_2(b^1\Sigma_g)$ plays an important role in the oxygen discharge

Thompson (1961) Proc. Royal Soc. A 262(1311) 519

- Recent global model study indicates there is a significant density of $O_2(b^1\Sigma_g)$ in the oxygen discharge
- The $O_2(b^1\Sigma_g)$ is mainly created through

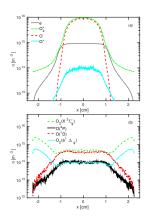
$$O_2(X^3\Sigma_g^-) + O(^1D) \longrightarrow O_2(b^1\Sigma_g) + O(^3P)$$

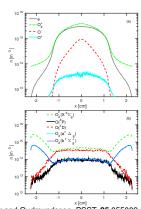


Toneli et al., J. Phys. D, 48 325202 (2015)



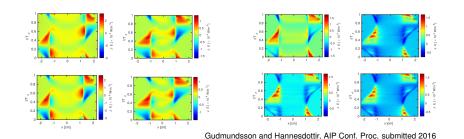






Hannesdottir and Gudmundsson, PSST, **25** 055002 (2016)

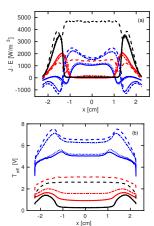
■ The density profiles of charged particles and fast neutrals comparing including $\gamma_{\text{see}}(E)$ and $O_2(a^1\Delta_g)$ (left) and $\gamma_{\text{see}}(E)$, $O_2(a^1\Delta_g)$ and $O_2(b^1\Sigma_g)$ (right) at 50 mTorr



- The spatio-temporal electron heating at 10 mTorr (left) and 50 mTorr (right)
- The four cases explored are:
 - (a) detachment neither by $O_2(a^1\Delta_g)$ nor $O_2(b^1\Sigma_g^+)$
 - \blacksquare (b) only detachment by $O_2(b^1\Sigma_g^+)$ included
 - \blacksquare (c) only detachment by $O_2(a^1\Delta_g)$ included
 - \blacksquare (d) both detachment by $O_2(a^1\Delta_g)$ and $O_2(b^1\Sigma_g^+)$ included

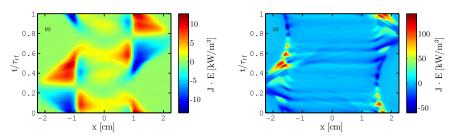


- The number of cold electrons increases as $O_2(b^1\Sigma_g)$ is added to the discharge model
- The electron heating in the bulk drops to zero
- The effective electron temperature profile changes significantly when detachment by singlet metastables is added to the reaction set
- 10 mTorr, 50 mTorr and 200 mTorr





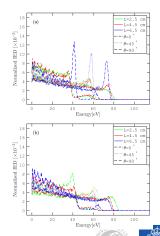




■ The spatio-temporal electron heating at 10 mTorr with a gap separation of 4.5 cm driven by a 222 V voltage source at 13.56 MHz (left) and 27.12 MHz (right)

Capacitively Coupled Oxygen Discharge Dual Frequency at 13.56 MHz + 27.12 MHz

- The ion energy distribution (IED)
- A dual frequency (13.56 MHz + 27.12 MHz) parallel plate capacitively coupled oxygen discharge at 75 mTorr
 - (a) Detachment by O₂(a¹Δ_g) and O₂(b¹Σ_g) included
 - (b) Detachment by O₂(a¹Δ_g) and O₂(b¹Σ_g) excluded
- For a larger discharge gap (*L* = 6.5 cm), the peak in the IED is much more apparent in the case where the full reaction set is used in the discharge model



Comparison to experimental findings:

o
$$\gamma_{\text{see}} = 0.0$$
,
4.4 % $O_2(a^1 \Delta_g)$

$$+ \gamma_{\text{see}} = 0.0,$$

4.4 % O₂(a¹ Δ_g) and 4.4 % O₂(b¹ Σ_g)

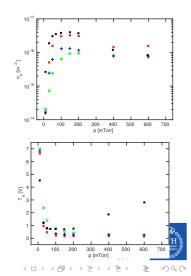
$$\mathbf{x} \ \gamma_{\text{see}} = \gamma_{\text{see}}(E),$$

4.4 % $O_2(\mathbf{a}^1 \Delta_g)$ and 4.4 % $O_2(\mathbf{b}^1 \Sigma_g)$

Experimental findings by Kechkar

(S. Kechkar, Ph.D. Thesis, Dublin City University, January 2015)

Hannesdottir and Gudmundsson (2016) PSST 25 055002



Summary

- We demonstrated particle-in-cell/Monte Carlo collision simulation of a capcacitively coupled disharge
- In an oxygen discharge at low pressure the EEPF is convex and develops a concave shape or becomes bi-Maxwellian as the pressure is increased
- These results contradict what is commonly found for the capacitively coupled argon discharge where the EEPF evolves from being concave at low pressure to being convex at high pressure





C. Overall Summary





Overall Summary

- A global (volume averaged) model can be used to understand the plasma chemistry
 - Which particles are important
 - Which reactions are important
 - How do the plasma parameters scale with the control parameters – power, pressure, discharge dimensions
- Particle-in-cell/Monte Carlo collision simulations can be used to explore the plasma kinetics
 - To find the electron energy distribution function
 - To find the ion energy distribution (IED) and the ion angular distribution (IAD)





Remaining questions

- There is clearly a need for basic experimental work on capacitively coupled oxygen discharges
 - Langmuir probe measurements, EEDF, *T*_{eff}, *n*_e versus pressure and applied voltage
 - Electronegativity for various pressures and applied voltage
- There are no cross sections available for electron impact excitation, dissociation or ionization from the singlet metastable states $O_2(a^1\Delta_g)$ and $O_2(b^1\Sigma_g)$
- The rate coefficient for detachment by $O_2(b^1\Sigma_g)$ has to be reevaluated





Remaining questions

- What is the role of electron impact dissociative attachment from vibrationally excited molecular states?
- What is the role of the metastable Herzberg states $O_2(A^3\Sigma_u^+, A'^3\Delta_u, c^1\Sigma_u^-)$?
- Wall recombination of O(³P) and in particular for O(¹D) as well as the wall quenching coefficient remain poorely known
- Measurements of the wall quenching coefficient of O₂(b¹Σ_a) are needed





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http://langmuir.raunvis.hi.is/~tumi/ranns.html

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