

Vlasov simulation of auroral processes

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Simulation model

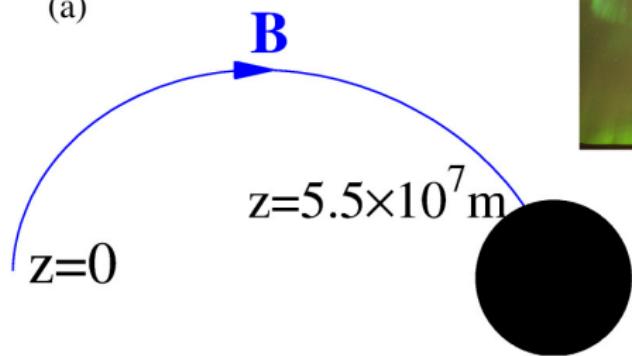
The distribution function is $f_s(z, v_z, \mu)$ and the Vlasov equation

$$\frac{\partial f_s}{\partial t} + v_z \frac{\partial f_s}{\partial z} + \frac{1}{m_s} \left(q_s E - \mu \frac{\partial B}{\partial z} + m_s a_g \right) \frac{\partial f_s}{\partial v_z} = 0$$



$$\mu = \frac{m_s v_{\perp}^2}{2B}$$

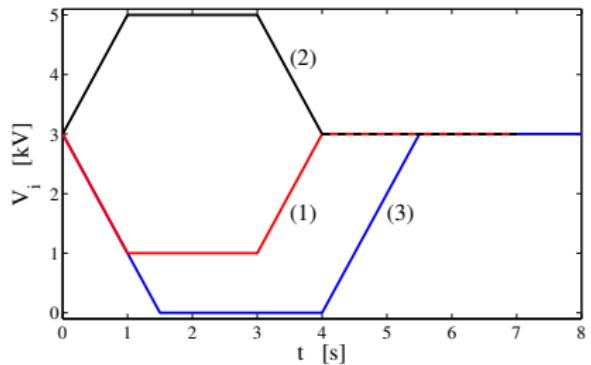
is a constant of motion.
The electric field is given by



$$\frac{d}{dz} \left(\frac{B_S}{B} E \right) = \frac{\rho_I}{S\epsilon}$$

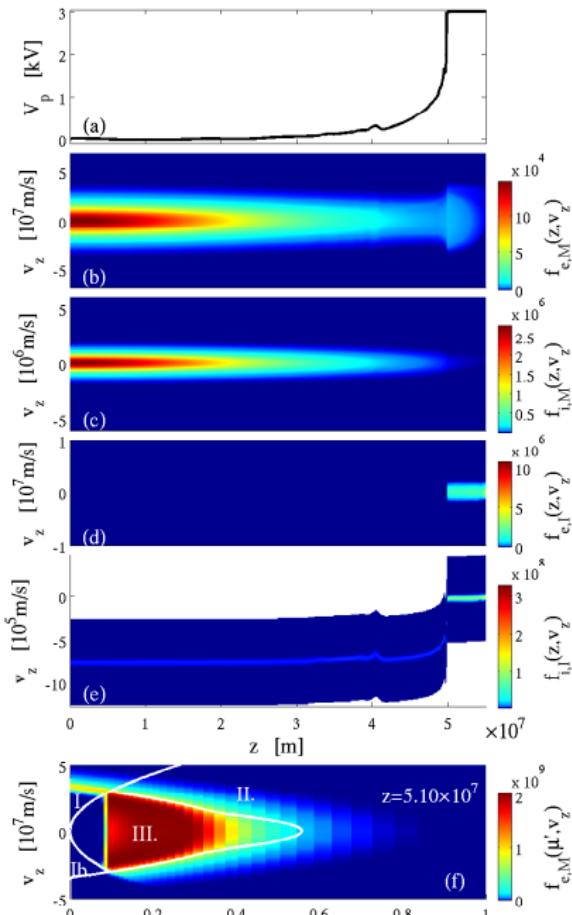
Computer code published
by Gunell et al. (2013a)

Changing the acceleration voltage during the simulation

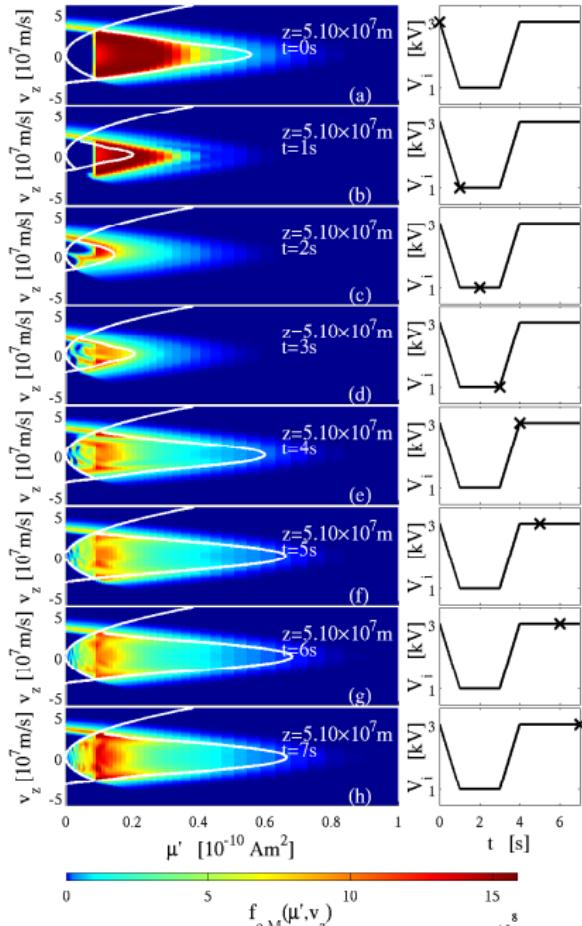


Above: Prescribed total voltage as a function of time.

Right: Initial state, common to all three experiments.

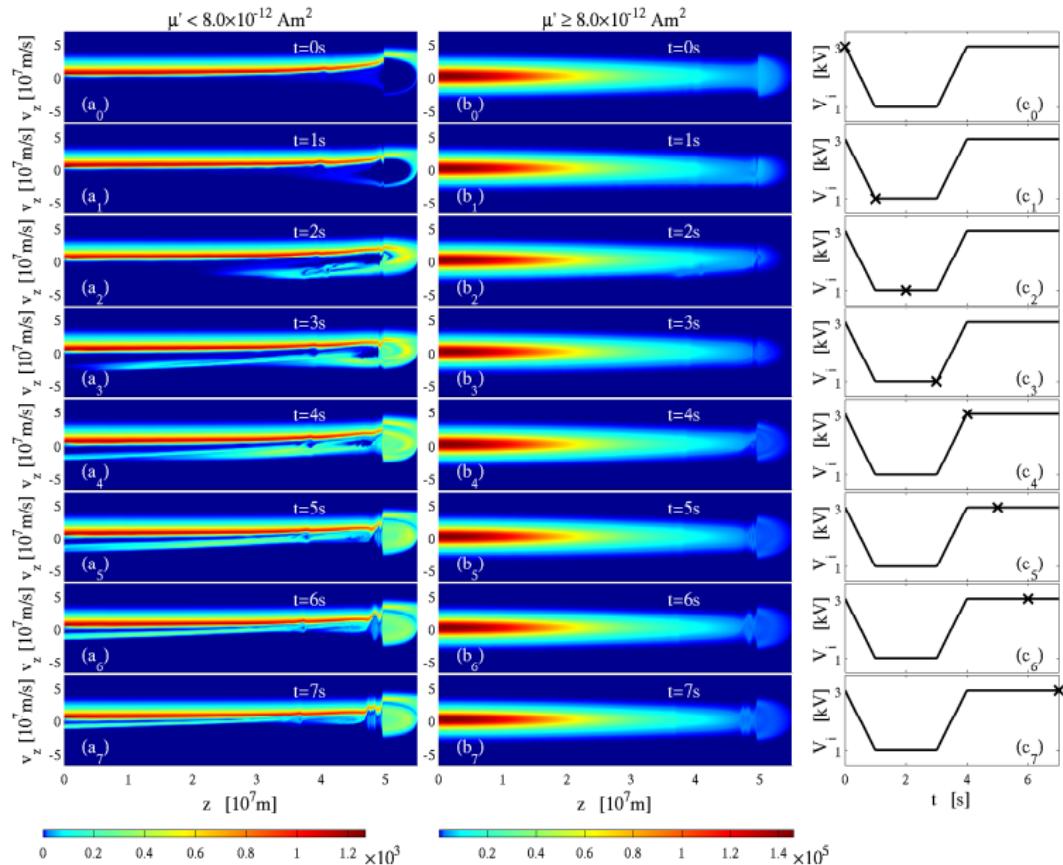


Experiment 1: $f(\mu, v_z)$ at $z = 5.1 \times 10^7$ m

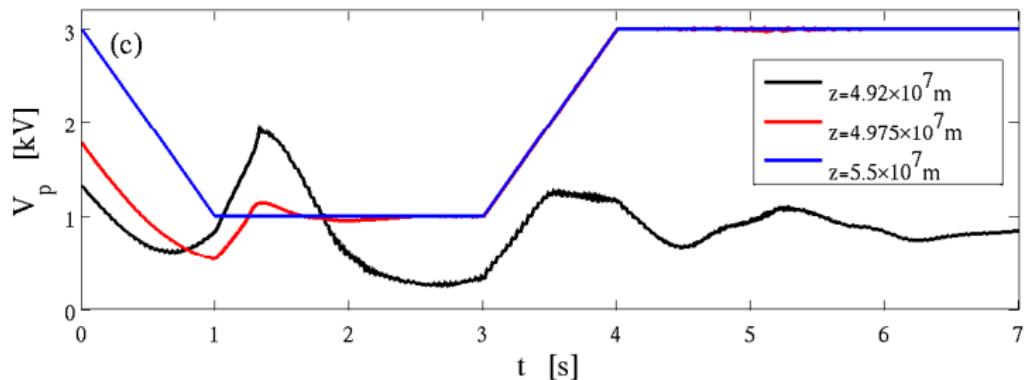
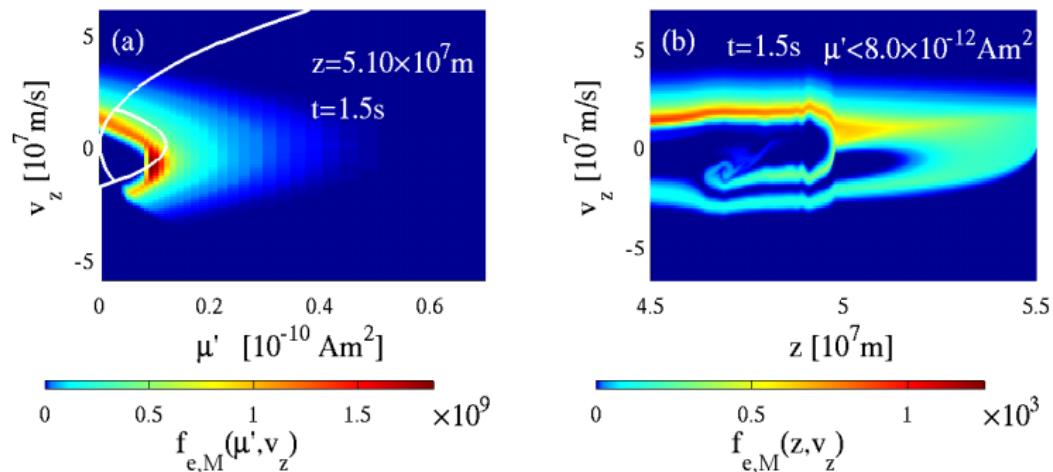


- ▶ Trapped electrons are released.
- ▶ New electrons get trapped.
- ▶ The distribution in the μ direction changes.

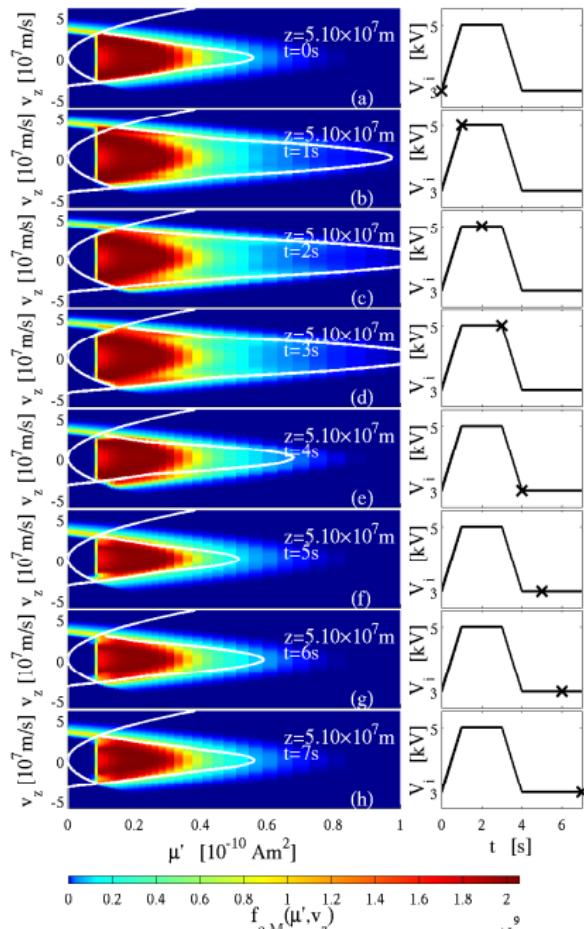
Experiment 1: $f(z, v_z)$



Experiment 1: fluctuations

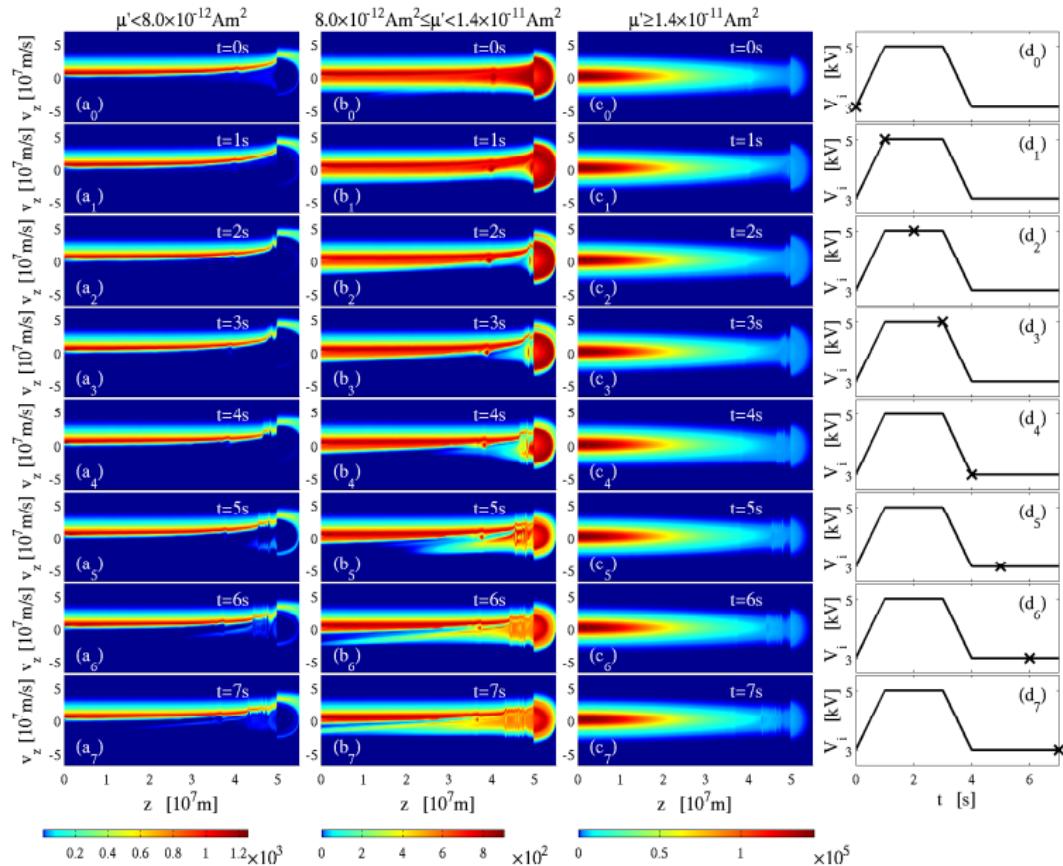


Experiment 2: $f(\mu, v_z)$ at $z = 5.1 \times 10^7$ m

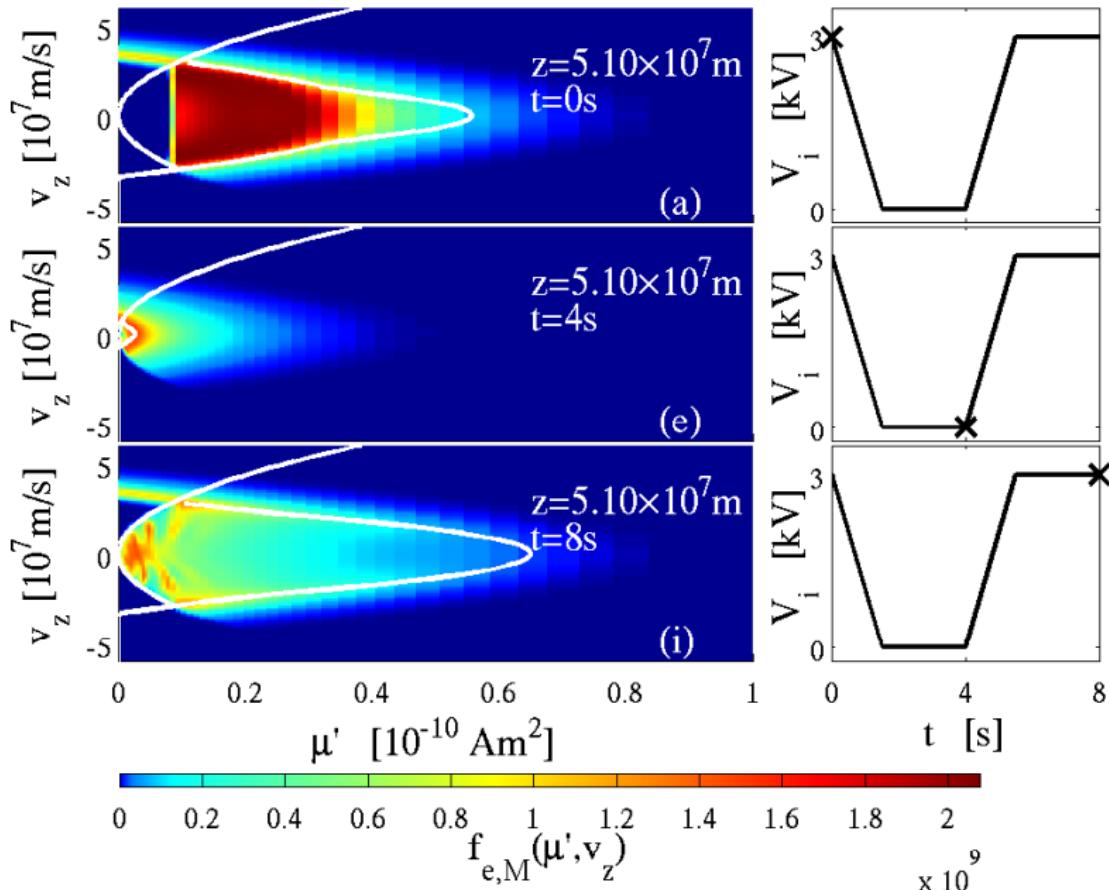


- ▶ New electrons are trapped at large μ values.
- ▶ Then these are lost again.
- ▶ The final state resembles the initial, although there is new structure at intermediate μ values.

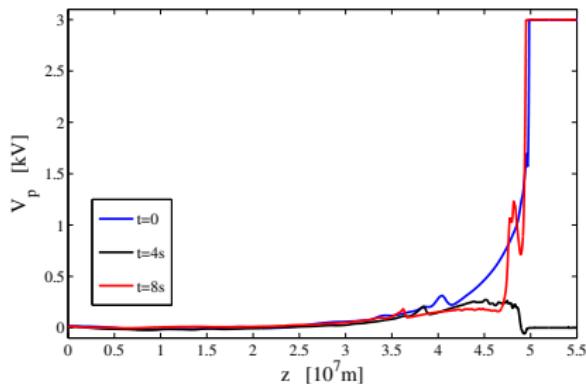
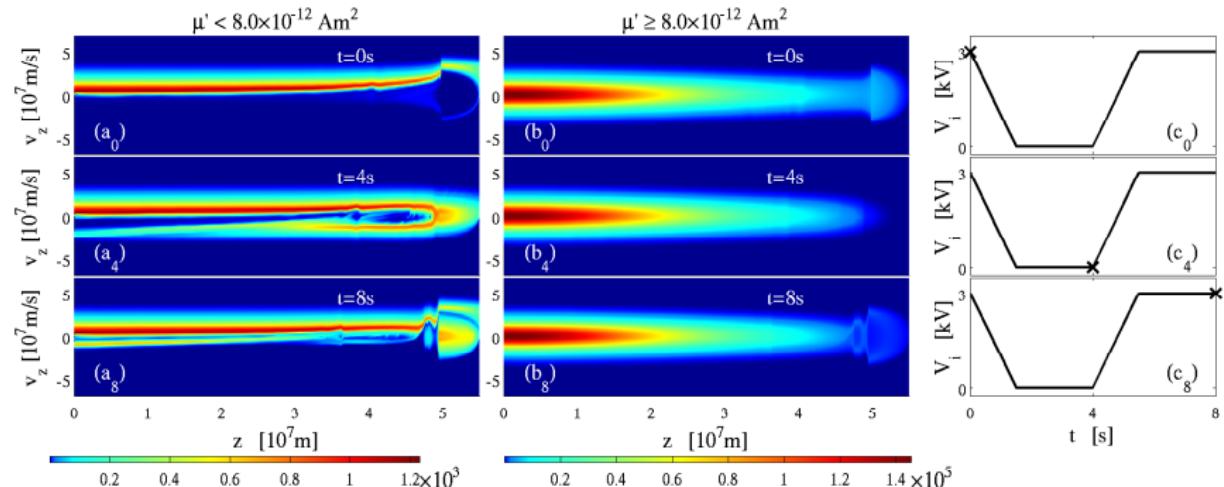
Experiment 2: $f(z, v_z)$



Experiment 3: $f(\mu, v_z)$ at $z = 5.1 \times 10^7$ m

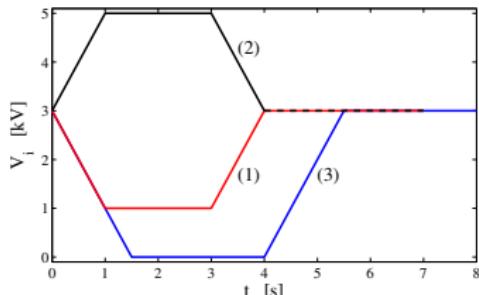
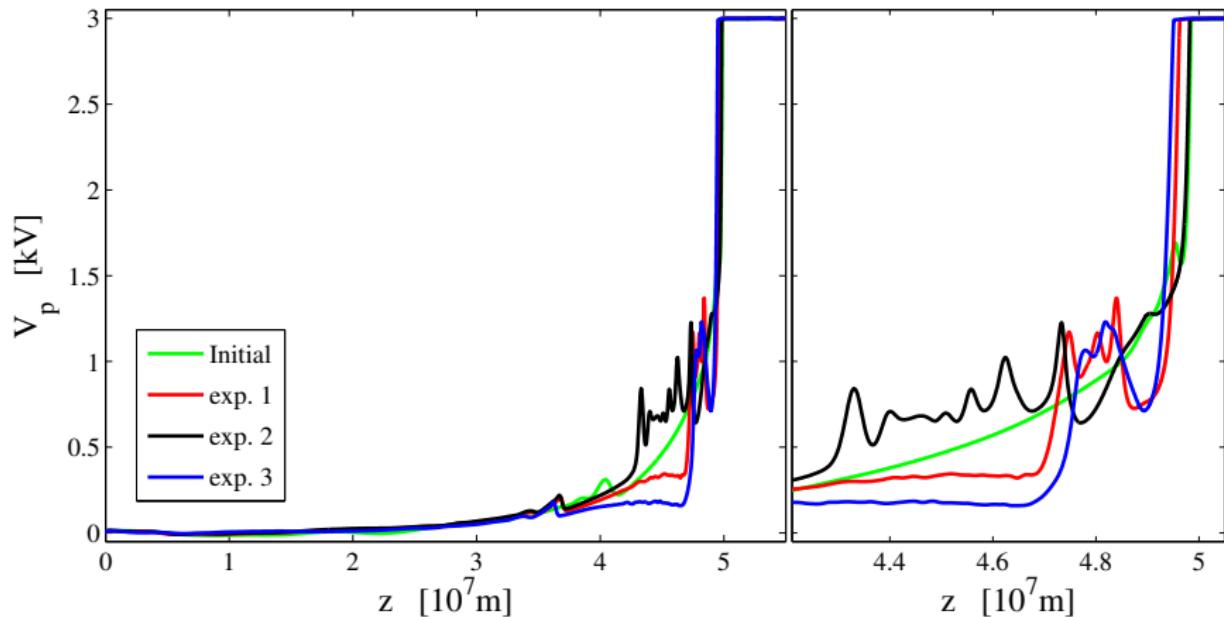


Experiment 3: $f(z, v_z)$



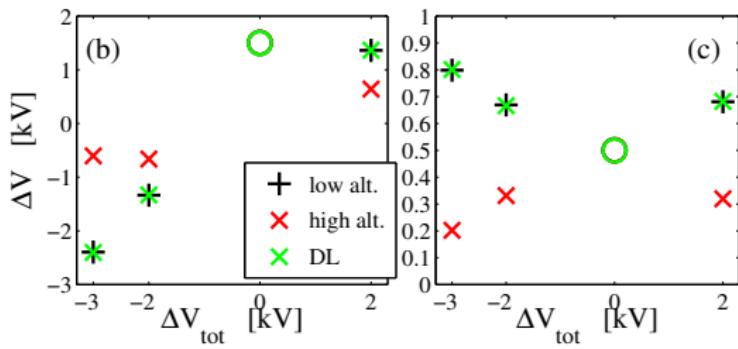
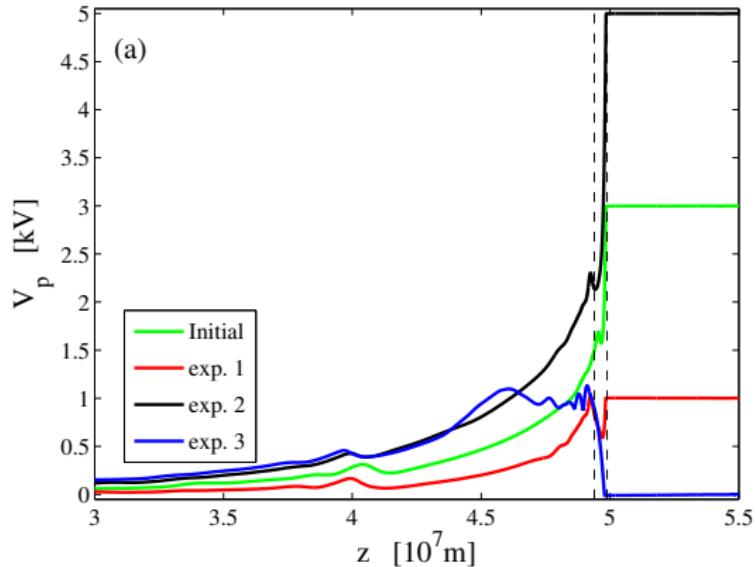
- ▶ The double layer changes polarity when the voltage is low.
- ▶ At $t = 8\text{s}$, a new potential peak traps some electrons.

Final state

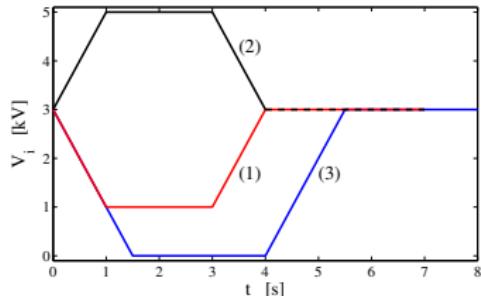


- ▶ The potential structure changes on the low potential side of the double layer.
- ▶ The double layer moved for $\Delta V < 0$.

Where does ΔV go?

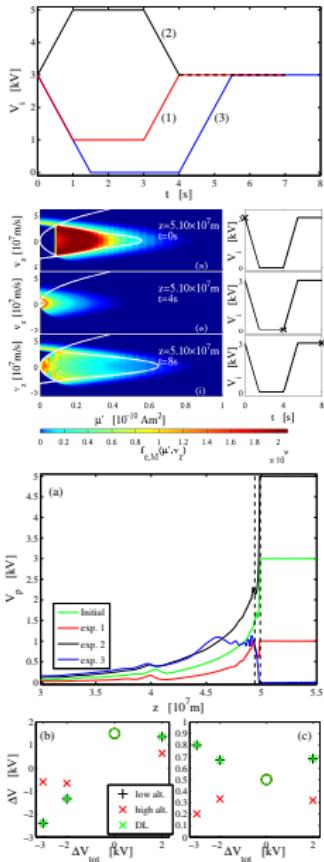


Most of ΔV is assumed by the double layer, c.f. observations by Forsyth et al. (2012).



Summary

- ▶ Trapping and release processes can be followed in simulations.
- ▶ The history of the acceleration voltage affects the distribution function of the trapped electrons.
- ▶ Most of the change in the acceleration voltage is assumed by the double layer.
- ▶ For a quickly decreasing voltage the double layer changes polarity.
- ▶ Phase space holes are created in these numerical experiments.
- ▶ The double layer position exhibits hysteresis phenomena.



Bibliography

- Forsyth, C., Fazakerley, A. N., Walsh, A. P., Watt, C. E. J., Garza, K. J., Owen, C. J., Constantinescu, D., Dandouras, I., Fornacon, K.-H., Lucek, E., Marklund, G. T., Sadeghi, S. S., Khotyaintsev, Y., Masson, A., Doss, N., Dec. 2012. Temporal evolution and electric potential structure of the auroral acceleration region from multispacecraft measurements. *Journal of Geophysical Research (Space Physics)* 117 (A16), 12203.
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- Gunell, H., De Keyser, J., Mann, I., 2013b. Numerical and laboratory simulations of auroral acceleration. *Physics of Plasmas* 20, 102901.
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Vlasov simulations – technical details – ϵ_r

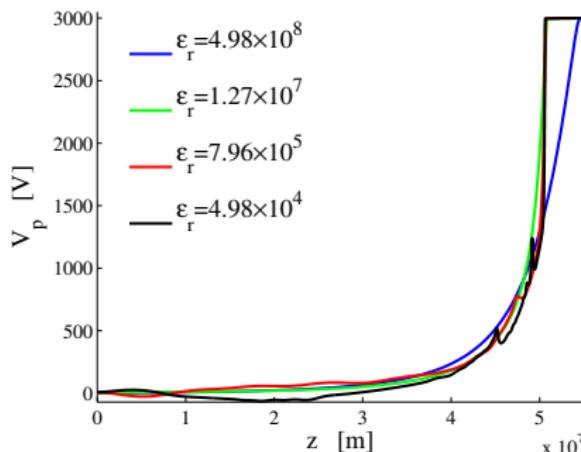
Computational advantages

$\omega_p \sim 1/\sqrt{\epsilon_r} \Rightarrow$ lets us use a longer time step

$\lambda_D \sim \sqrt{\epsilon_r} \Rightarrow$ lets us use a coarser grid

Disadvantage

Suppression of small scale and high frequency phenomena



	Magneto-sphere	Iono-sphere
z	0	$5.5 \cdot 10^7$ m
B	$0.086 \mu\text{T}$	$56 \mu\text{T}$
$k_B T_e$	500 eV	1 eV
$k_B T_{H^+}$	2500 eV	1 eV
n	$3 \cdot 10^5 \text{ m}^{-3}$	$1 \cdot 10^9 \text{ m}^{-3}$