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## Modelling radical production during the TORCH campaign

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The Tropospheric ORganic CHemistry experiment (TORCH) took place during the unusually hot July and August of 2003 at Writtle College, a site 2 miles west of Chelmsford in Essex and 20 miles North East of London. The concentrations of the hydroxyl radical OH, hydroperoxy radical HO $_2$  and alkyl peroxy radicals RO $_2$ , were measured during the campaign. Predictions were also made using a zero-dimensional box model using the Master Chemical Mechanism (v3.1) constrained with the concentrations of relatively long-lived species. New model parameterisations include a heterogeneous loss rate for HO $_2$  and a production mechanism for HONO. The model results were compared to radical measurements made by FAGE and PERCA. The modelled OH results were in good agreement with the measurements. The dominant OH source was from the reaction of J(O $^1$ D) and H $_2$ O during the day, with the O $_3$  -alkenes reaction producing most OH during the night.

The 15-minute average diurnal time series for modelled and measured  $HO_2$  showed agreement in shape and magnitude. The relationship between the modelled  $HO_2/OH$  ratio and measured NO concentrations was stronger than that of the measured  $HO_2/OH$  ratio and NO. On average NO concentrations were 2.8 ppb during the campaign, with a peak of 44.7 ppb on July  $28^{th}$ . The modelled peroxy radical group  $\Sigma[HO_2+RO_2]$  was under-predicted when compared with their respective measurements from PERCA. The under-prediction was always greater during the daylight periods of the day, and especially at noon. With such good agreement between modelled and measured  $HO_2$  concentrations, these results suggest that there are too few sources of  $RO_2$  radicals in the model. These results are discussed in context with modelling studies conducted for the urban PUMA campaign.