

**Table S1:** Values of the parameters used to calculate the modelled emissions with L+T or T algorithms, standardised emissions factors for *L+T* algorithm ( $EF_{L+T}$ ), specific standardised emissions factors for *T* algorithm ( $EF_T$ ) and experimental coefficient  $\beta$ . Means  $\pm$  se, n = 5.

**Table S1:**

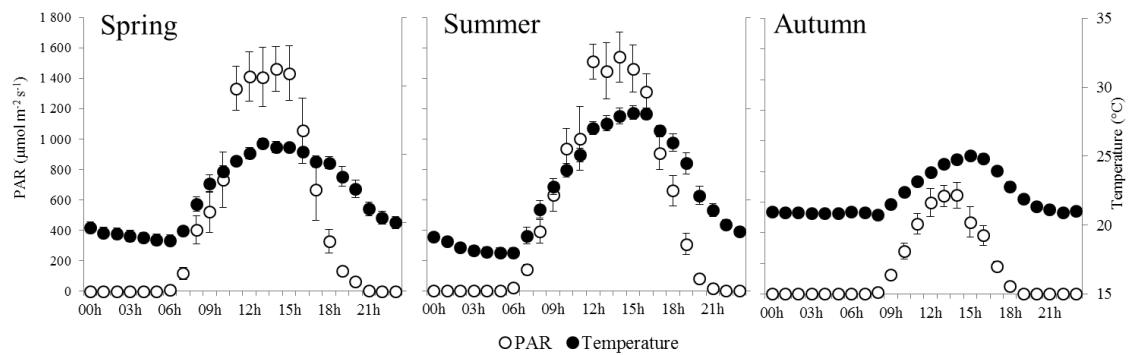
Compounds	Treatment	Spring			Summer			Autumn		
		$EF_{L+T}$	$EF_T$	$\beta$	$EF_{L+T}$	$EF_T$	$\beta$	$EF_{L+T}$	$EF_T$	$\beta$
<b>Isoprene</b>	<b>ND</b>	28.5 $\pm$ 4.6			118.0 $\pm$ 8.4			6.4 $\pm$ 1.1		
	<b>AD</b>	17.8 $\pm$ 2.4			84.8 $\pm$ 9.0			12.0 $\pm$ 2.8		
<b>MACR+MVK+ISOPOOH</b>	<b>ND</b>	0.2 $\pm$ 0.03	0.7 $\pm$ 0.15	0.5 $\pm$ 0.01	0.3 $\pm$ 0.02	0.4 $\pm$ 0.1	0.5 $\pm$ 0.1	0.1 $\pm$ 0.01	1.6 $\pm$ 0.3	0.7 $\pm$ 0.1
	<b>AD</b>	0.1 $\pm$ 0.01	0.9 $\pm$ 0.28	0.6 $\pm$ 0.04	0.2 $\pm$ 0.03	0.2 $\pm$ 0.04	0.6 $\pm$ 0.1	0.1 $\pm$ 0.02	5.0 $\pm$ 1.6	0.9 $\pm$ 0.03
<b>Methanol</b>	<b>ND</b>	1.0 $\pm$ 0.2	2.6 $\pm$ 0.8	0.3 $\pm$ 0.1	0.7 $\pm$ 0.04	0.9 $\pm$ 0.04	0.3 $\pm$ 0.04	0.3 $\pm$ 0.1	1.1 $\pm$ 0.3	0.4 $\pm$ 0.1
	<b>AD</b>	0.8 $\pm$ 0.1	2.2 $\pm$ 0.2	0.3 $\pm$ 0.04	0.5 $\pm$ 0.1	0.6 $\pm$ 0.04	0.3 $\pm$ 0.1	0.4 $\pm$ 0.1	1.4 $\pm$ 0.4	0.4 $\pm$ 0.04
<b>Acetone</b>	<b>ND</b>	0.6 $\pm$ 0.2	1.8 $\pm$ 0.5	0.4 $\pm$ 0.01	0.9 $\pm$ 0.1	1.1 $\pm$ 0.2	0.4 $\pm$ 0.1	0.6 $\pm$ 0.2	2.4 $\pm$ 0.6	0.4 $\pm$ 0.03
	<b>AD</b>	0.5 $\pm$ 0.1	2.1 $\pm$ 0.4	0.5 $\pm$ 0.02	0.4 $\pm$ 0.03	0.5 $\pm$ 0.1	0.3 $\pm$ 0.1	0.8 $\pm$ 0.3	4.3 $\pm$ 1.8	0.5 $\pm$ 0.1
<b>Formaldehyde</b>	<b>ND</b>	0.3 $\pm$ 0.1	0.8 $\pm$ 0.2	0.4 $\pm$ 0.02	0.3 $\pm$ 0.03	0.4 $\pm$ 0.1	0.4 $\pm$ 0.1	0.4 $\pm$ 0.1	1.6 $\pm$ 0.4	0.4 $\pm$ 0.04
	<b>AD</b>	0.2 $\pm$ 0.02	1.3 $\pm$ 0.2	0.5 $\pm$ 0.03	0.2 $\pm$ 0.03	0.2 $\pm$ 0.1	0.4 $\pm$ 0.2	0.6 $\pm$ 0.1	2.7 $\pm$ 0.7	0.5 $\pm$ 0.02
<b>Acetaldehyde</b>	<b>ND</b>	2.4 $\pm$ 0.7	9.4 $\pm$ 2.7	0.5 $\pm$ 0.03	1.7 $\pm$ 0.4	1.6 $\pm$ 0.3	0.4 $\pm$ 0.1	2.5 $\pm$ 0.6	34.0 $\pm$ 3.1	0.7 $\pm$ 0.02
	<b>AD</b>	2.1 $\pm$ 0.9	7.9 $\pm$ 4.2	0.5 $\pm$ 0.1	0.9 $\pm$ 0.1	0.9 $\pm$ 0.1	0.5 $\pm$ 0.1	2.5 $\pm$ 0.7	37.8 $\pm$ 4.3	0.7 $\pm$ 0.1

**Figure S1:** Diurnal pattern of photosynthetic active radiations (PAR) and temperatures in spring, summer and autumn. Values are mean  $\pm$  SE, n=5

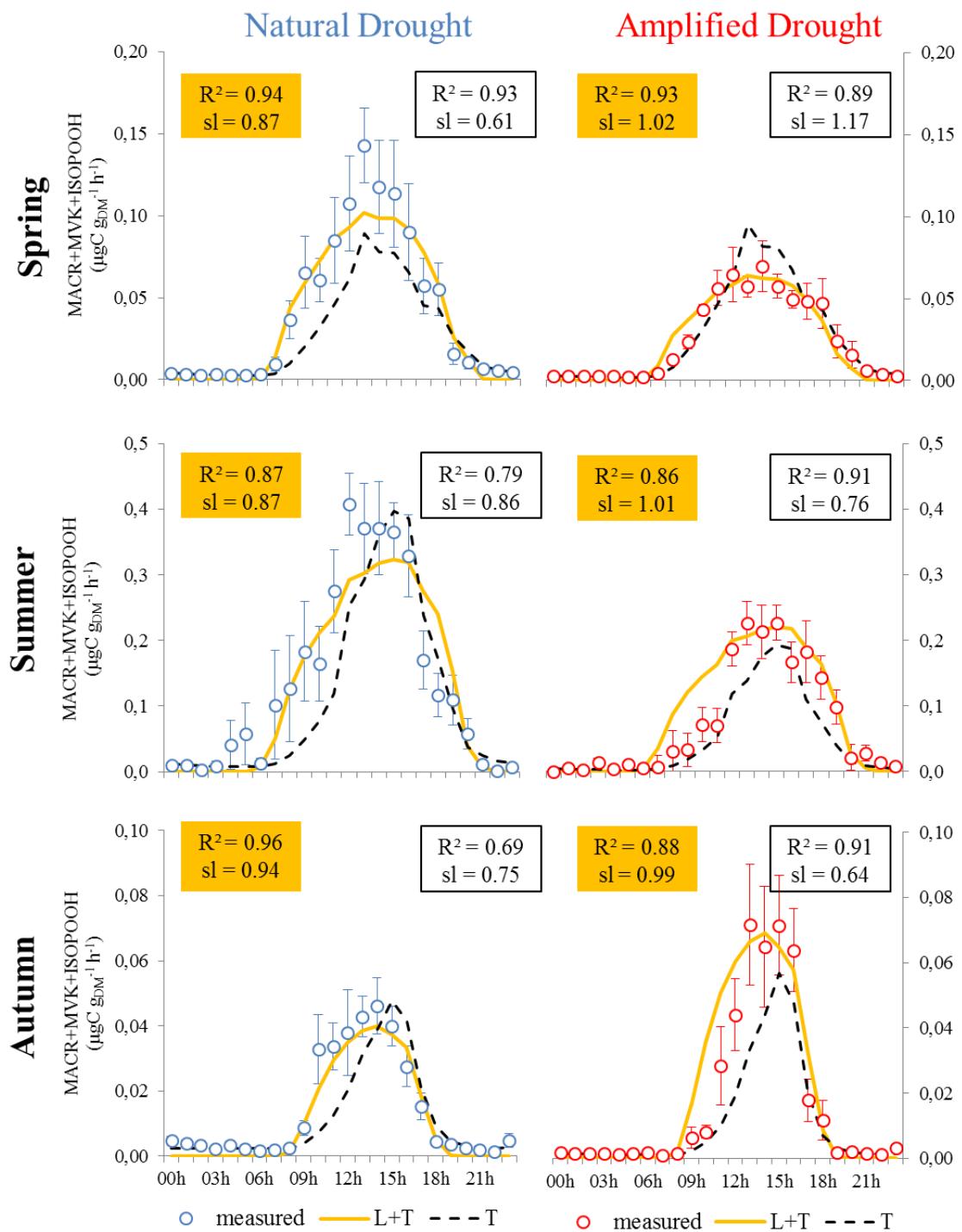
**Figure S2:** Diurnal pattern of MACR+MVK+ISOPPOOH emissions rates, where points represent measured emissions, the yellow line correspond to modelled emissions rates according to the  $L+T$  algorithm ( $ER_{L+T}$ ) and dotted line is modelled emissions rates according to  $T$  algorithm ( $ER_T$ ). Values are mean  $\pm$  SE, n=5.  $R^2$  and slope (sl) of correlations between measured and modelled emissions were presented in the yellow frame for  $L+T$  and in the white frame for  $T$ . Correlations were obtained without forcing data through the origin.

**Figure S3:** Diurnal pattern of acetone emissions rates where points represent measured emissions, yellow line correspond to modelled emissions rates according to the  $L+T$  algorithm ( $ER_{L+T}$ ) and dotted line is modelled emissions rates according to  $T$  algorithm ( $ER_T$ ). Values are mean  $\pm$  SE, n=5.  $R^2$  and slope (sl) of correlations between measured and modelled emissions were presented in the yellow frame for  $L+T$  and in the white frame for  $T$ . Correlations were obtained without forcing data through the origin.

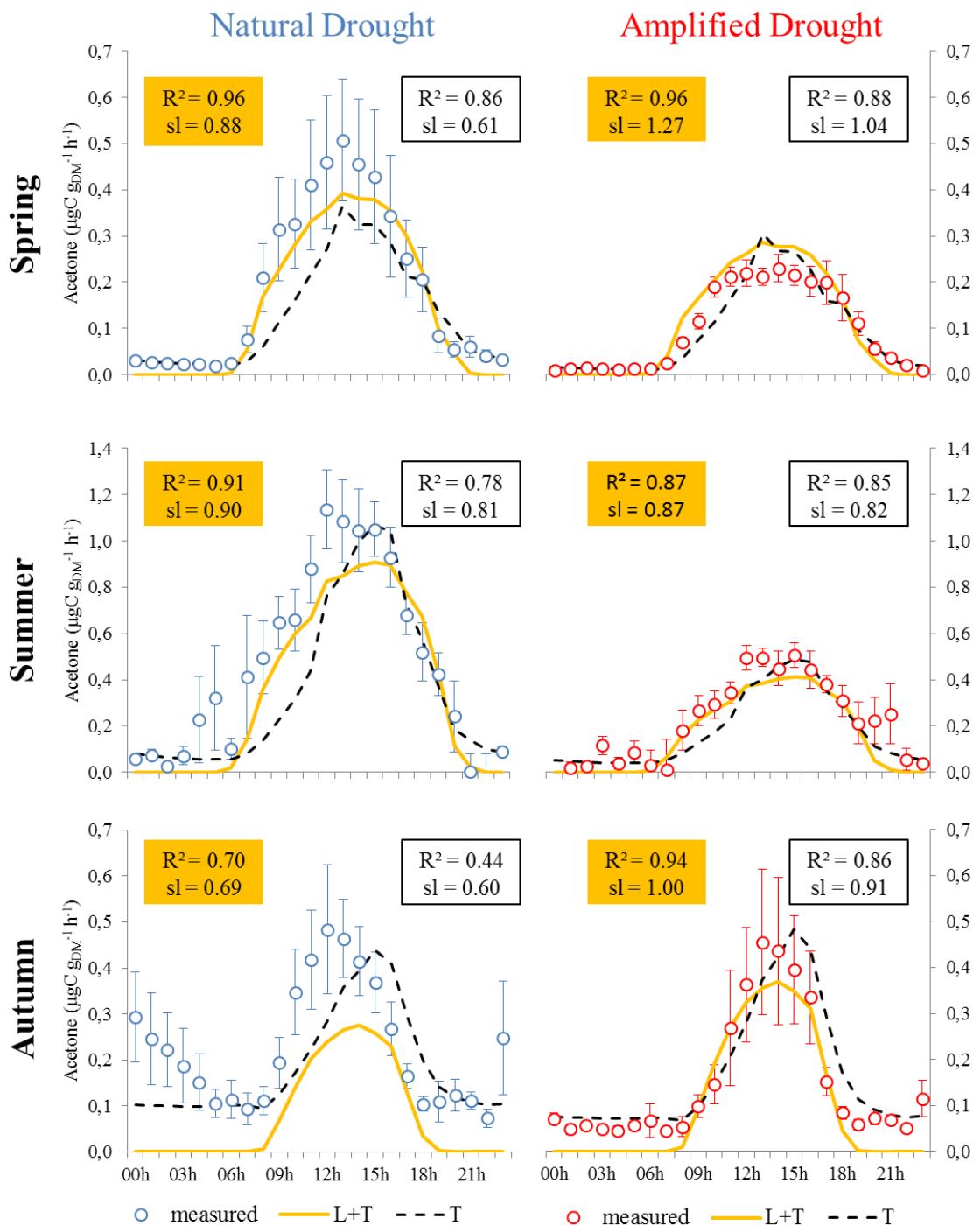
**Figure S4:** Diurnal pattern of formaldehyde emissions rates where points represent measured emissions, yellow line correspond to modelled emissions rates according to the  $L+T$  algorithm ( $ER_{L+T}$ ) and dotted line is modelled emissions rates according to  $T$  algorithm ( $ER_T$ ). Values are mean  $\pm$  SE, n=5.  $R^2$  and slope (sl) of correlations between measured and modelled emissions were presented in yellow frame for  $L+T$  and in white frame for  $T$ . Correlations were obtained without forcing data through the origin.



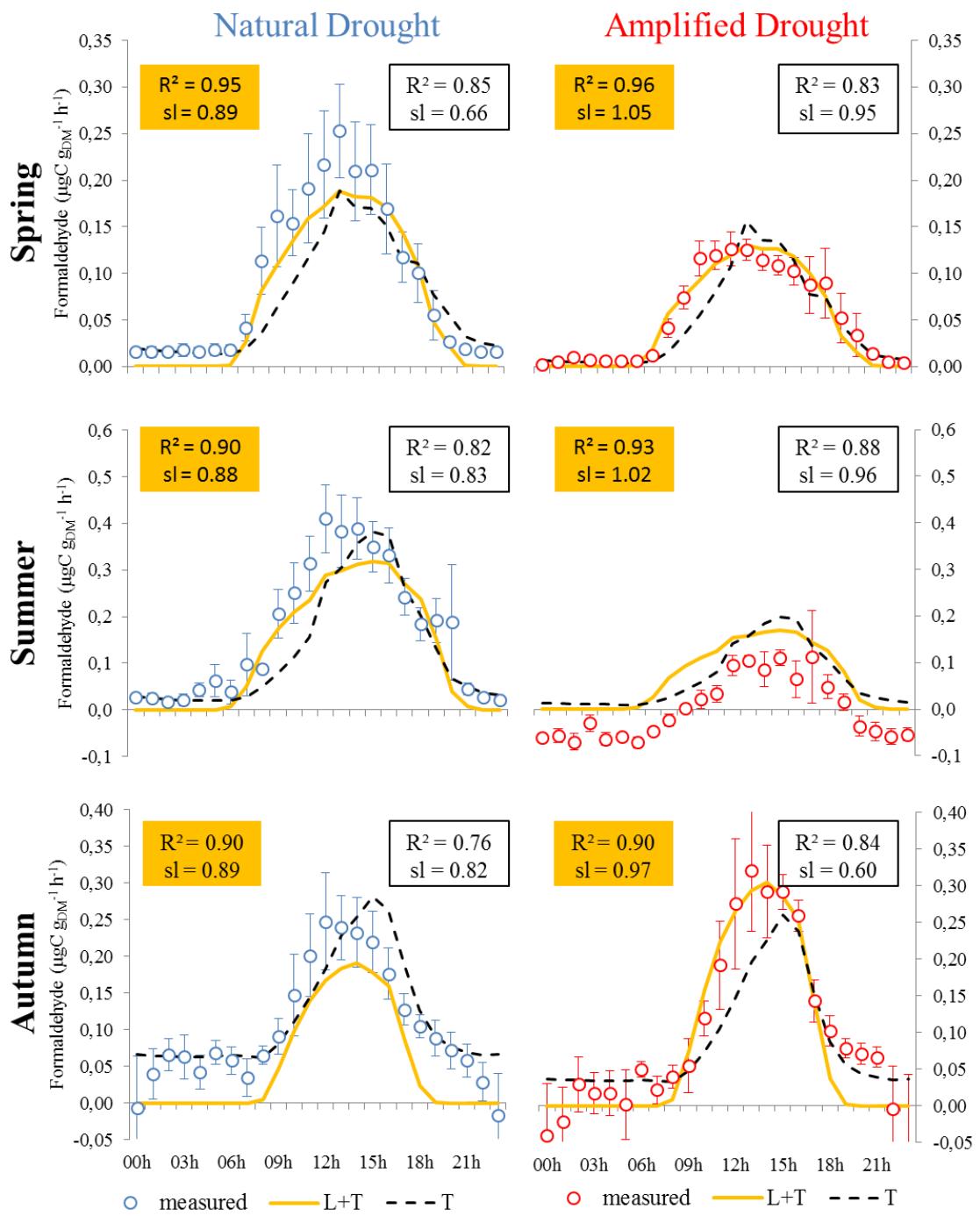
**Figure S1:**



**Figure S2:**



**Figure S3:**



**Figure S4:**

