

Identification of volatile organic compounds emitted by Sitka spruce and determination of their emission pathways and fluxes

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Supplementary Information

Set-up

Individual cylindrical Teflon enclosures were made for each spruce tree, using clear 0.1 mm thick Teflon FEP foil (Ortega and Helmig, 2008). The enclosures were designed to encompass all except the bottom few branches of the trees, and to minimise contact between the contained branches and the interior wall of the enclosure (Ortega and Helmig, 2008; Janson et al., 1999). The tops of the enclosures were fitted with a ¼” inlet port and a ½” outlet port. A length of ¼” Teflon tubing extended 20 cm from the inlet port into the enclosure to ensure homogenisation (Niinemets et al., 2011). The enclosures were sealed around the trees a week before emissions sampling was commenced. A separate, Empty enclosure, was made for background measurements (Hayward et al., 2004) (Fig. S3). ¼



Figure S1 Three Sitka spruce trees used in this work: Spruce 1 (left), Spruce 2 (middle) and Spruce 3 (right).

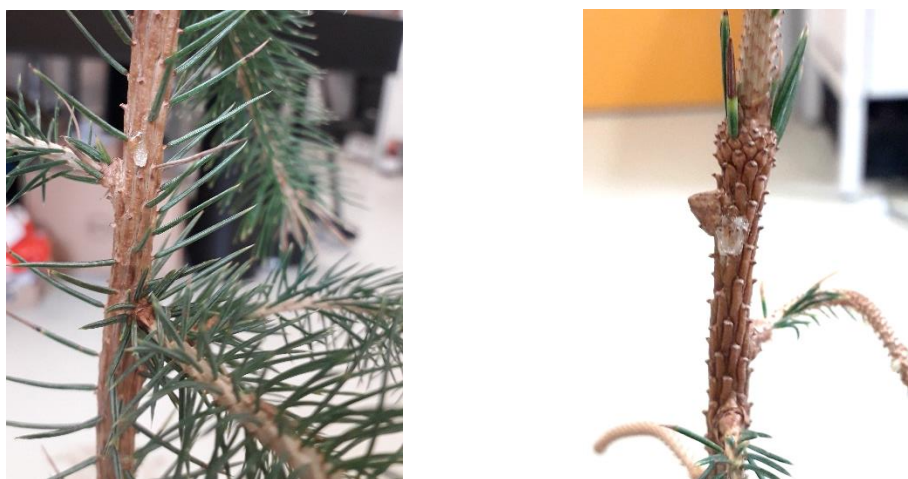


Figure S2 Resin deposits on Spruce 1 (left) and Spruce 3 (right).



Figure S3 Enclosure around Spruce 3 (left), and empty enclosure (right).

Plant growth cycles

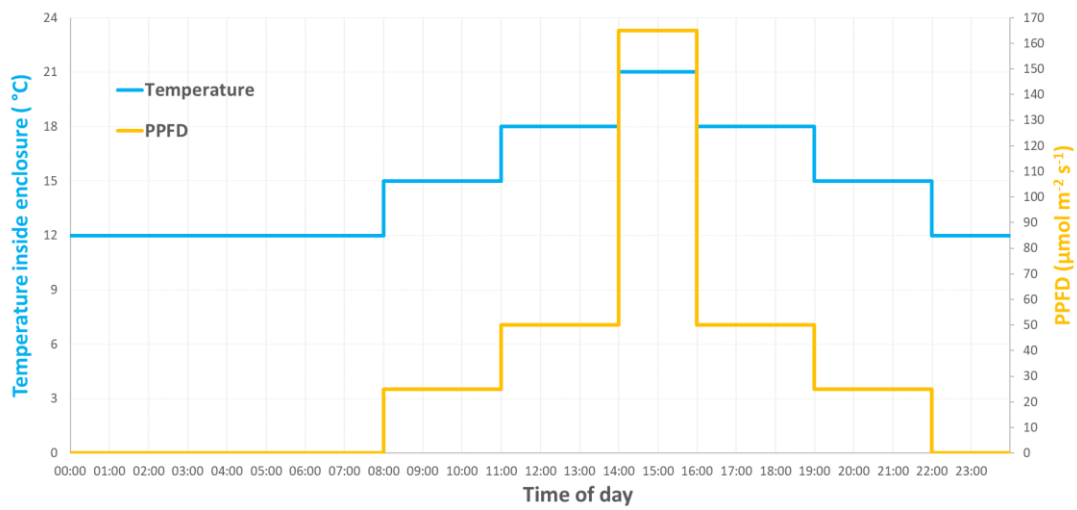


Figure S4 Temperature and PPFD variation during the Daily Cycle.

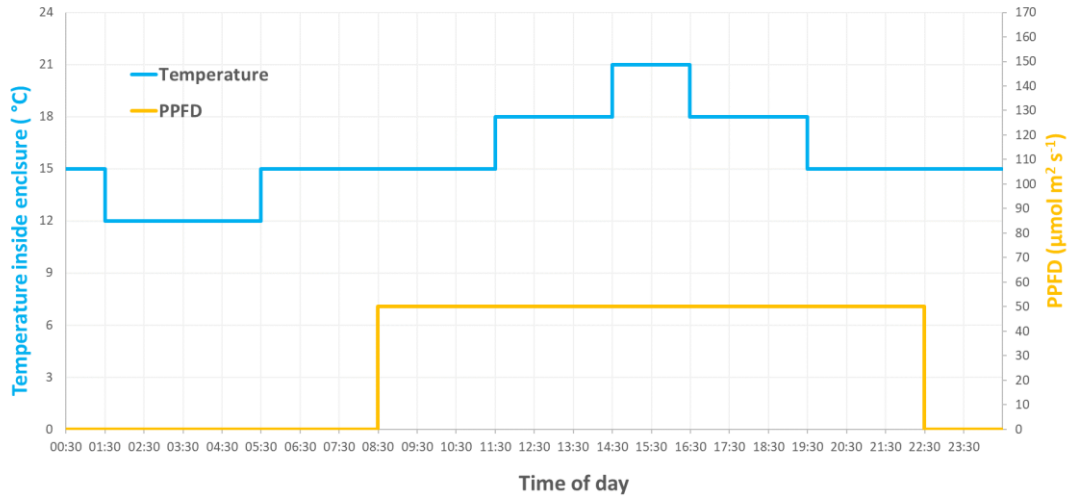


Figure S5 Temperature and PPFD variation during the Temperature Cycle.

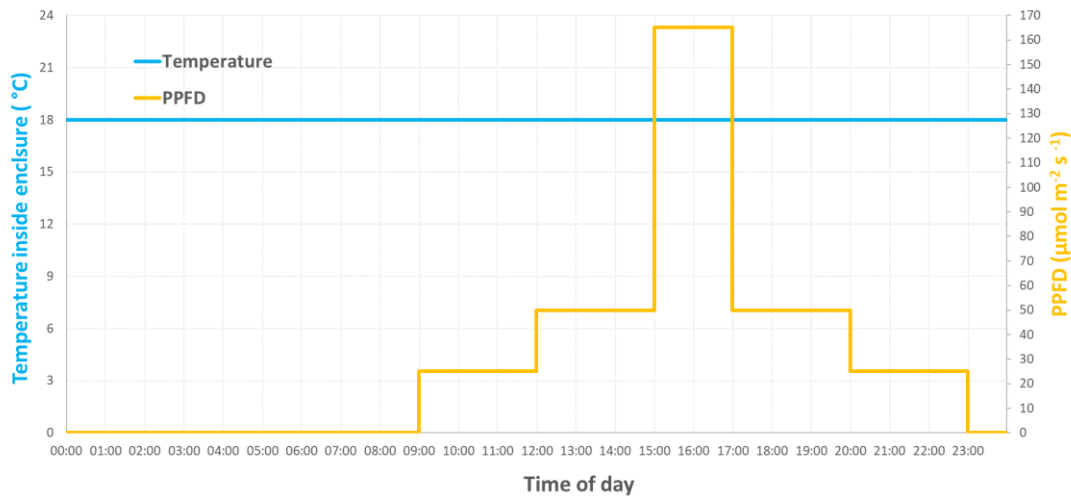


Figure S6 Temperature and PPFD variation during the Light Cycle.

Biomass measurements

Table S1 Dried biomass measurements for each spruce tree.

	Dried Needle Mass	Dried Needle and Branch Mass
	g	g
Spruce 1	8.03	18.50
Spruce 2	9.80	19.56
Spruce 3	3.77	11.33

BVOC emissions

Table S2 BVOC emissions from Spruce 1 and Spruce 2 detected with ToF-CIMS and TD-GC/MS.

Formula	MW g mol ⁻¹	Name	Spruce 1		Spruce 2	
			ToF-CIMS	TD-CG/MS	ToF-CIMS	TD-GC/MS
C ₃ H ₉ N	59.11		✓			
C ₂ H ₄ O ₂	60.05	Acetic acid				✓
C ₅ H ₈	68.12	Isoprene	✓	✓		✓
C ₄ H ₆ O	70.09	Methacrolein				✓
C ₃ H ₆ O ₂	74.08	Propanoic acid		✓		✓
C ₄ H ₁₀ O	74.12	1-Butanol				✓
C ₆ H ₈	80.13				✓	
C ₅ H ₆ O	82.10	2-Methyl-furan		✓		✓
C ₅ H ₈ O	84.12	<i>E</i> -3-Penten-2-one				✓
C ₅ H ₁₀ O	86.13	2-Methyl-3-buten-2-ol		✓		
		2-Pentanone		✓		✓
		Pentanal				✓
C ₄ H ₈ O ₂	88.11	Butanoic acid				✓
C ₃ H ₈ N ₂ O	88.11		✓			
C ₅ H ₁₂ O	88.15	3-Methyl-1-butanol		✓		✓
C ₇ H ₈	92.14		✓		✓	
C ₆ H ₆ O	94.11	Phenol				✓
C ₆ H ₈ O	96.13				✓*	
C ₆ H ₁₀ O	98.15	<i>E</i> -2-Hexenal			✓	✓
C ₅ H ₈ O ₂	100.13		✓			
C ₆ H ₁₂ O	100.16	Hexanal				✓
		2-Hexanone				✓
C ₇ H ₁₆	100.20	Heptane		✓		✓
C ₆ H ₁₄ O	102.18	1-Hexanol		✓		✓
C ₇ H ₈ O	108.14		✓		✓	
C ₈ H ₁₂	108.18		✓		✓	
C ₇ H ₁₀ O	110.15				✓	
C ₅ H ₁₀ NO ₂	117.15		✓			
C ₈ H ₈ O	120.15	Acetophenone				✓
C ₉ H ₁₂	120.19		✓		✓	
C ₇ H ₁₂ N ₂	124.18		✓			
C ₈ H ₁₂ O	124.18				✓	
C ₈ H ₁₆ O	128.21	1-Octen-3-ol				✓
C ₁₀ H ₁₀	130.19		✓			
C ₈ H ₁₈ O	130.23	2-ethyl-1-hexanol				✓
C ₁₀ H ₁₄	134.22	<i>o</i> -Cymene		✓		✓
C ₉ H ₈ O	132.16	(<i>E</i>)-Cinnamaldehyde	✓	✓	✓	✓
C ₆ H ₁₄ O ₃	134.18		✓			

C ₈ H ₈ O ₂	136.15	Methyl benzoate		✓		✓
C ₉ H ₁₂ O	136.19				✓	
C ₁₀ H ₁₆	136.24	Monoterpene	✓		✓*	
		Myrcene		✓		✓
		β-Phellandrene		✓		✓
		δ-Limonene		✓		✓
		α-Pinene		✓		✓
		Camphene		✓		✓
C ₈ H ₁₂ O ₂	140.18				✓	
C ₉ H ₁₀ O ₂	150.17				✓	
C ₃ H ₉ N	150.22		✓			
C ₇ H ₄ O ₄	152.10		✓			
C ₈ H ₈ O ₃	152.14	Methyl salicylate		✓	✓	✓
C ₁₀ H ₁₆ O	152.23	Piperitone	✓	✓	✓	✓
		Camphor	✓	✓	✓	✓
C ₁₀ H ₁₈ O	154.25	Eucalyptol		✓	✓	✓
C ₁₂ H ₁₆	160.25		✓			
C ₁₁ H ₁₄ O	162.23		✓			
C ₅ H ₁₀ O ₆	166.13		✓			
C ₁₀ H ₁₄ O ₂	166.22		✓		✓*	
C ₁₀ H ₁₆ O ₂	168.24		✓			
C ₁₀ H ₂₀ O ₂	172.26	Isopentyl isovalerate	✓	✓		✓
C ₁₂ H ₁₆ O	176.26				✓	
C ₁₂ H ₂₀ O	180.29		✓			
C ₁₀ H ₁₆ O ₃	184.23				✓	
C ₁₃ H ₁₂ O	184.24		✓			
C ₁₄ H ₁₆ O	200.28		✓		✓	
C ₁₀ H ₁₈ O ₄	202.25		✓		✓	
C ₁₅ H ₂₄	204.36	<i>E</i> -β-Farnesene			✓	✓
C ₁₀ H ₁₄ O ₅	214.22		✓			
C ₁₇ H ₂₆ O	246.39		✓		✓	
C ₂₀ H ₂₄	264.40		✓		✓	
C ₁₃ H ₃₀ O ₅	266.37		✓		✓	
C ₁₈ H ₃₄ O	266.46		✓		✓	
C ₂₀ H ₂₄ O ₂	296.40		✓		✓	

* Isomers with different temporal trends detected by ToF-CIMS as different ions. Counted as two distinct BVOCs. Monoterpenes are ordered according to decreasing concentration, which were inferred from TD-GC/MS chromatogram peak areas.

Chlorophyll fluorescence

Maximum photosystem II efficiency (F_v/F_m) was determined by measuring chlorophyll fluorescence. F_v/F_m reflects the ability of photosystem II to reduce its electron acceptors in the photosynthetic electron transport chain and it is used to infer plant health (Henriques, 2009). The chlorophyll fluorescence was measured on three days during the *Daily Cycle* (Table 2).

Table S3 Chlorophyll fluorescence measurements for Spruce 1 and Spruce 2 during the Daily Cycle. Each value is the average of three measurements taken from three branches on the same tree.

Date	PPFD $\mu\text{mol m}^{-2} \text{s}^{-1}$	Temperature $^{\circ}\text{C}$	Spruce 1		Spruce 2	
			F_v/F_m	STD	F_v/F_m	STD
04/11/2021	0	12	0.85	0.02	0.75	0.01
08/11/2021	0	12	0.79	0.02	0.83	0.03
11/11/2021	0	12	0.81	0.02	0.82	0.08

Chlorophyll fluorescence was measured in the dark period. The F_v/F_m values were similar for both Spruce 1 and Spruce 2 and was centred around 0.81 (unitless). However, for both trees values below 0.8 were recorded on one of the days, suggesting that there was some extent of photoinhibition due to stress. Lower F_v/F_m ratios indicate that all chlorophyll is not taking part in photosynthesis, which could be due to damage, and implies the plant is stressed (Maxwell and Johnson, 2000). It is unlikely the trees were suffering from stress, as the temperatures used in this study are representative of the natural environment, and the PPFD was an order of magnitude lower than that measured for the ambient atmosphere.

Temperature cycle

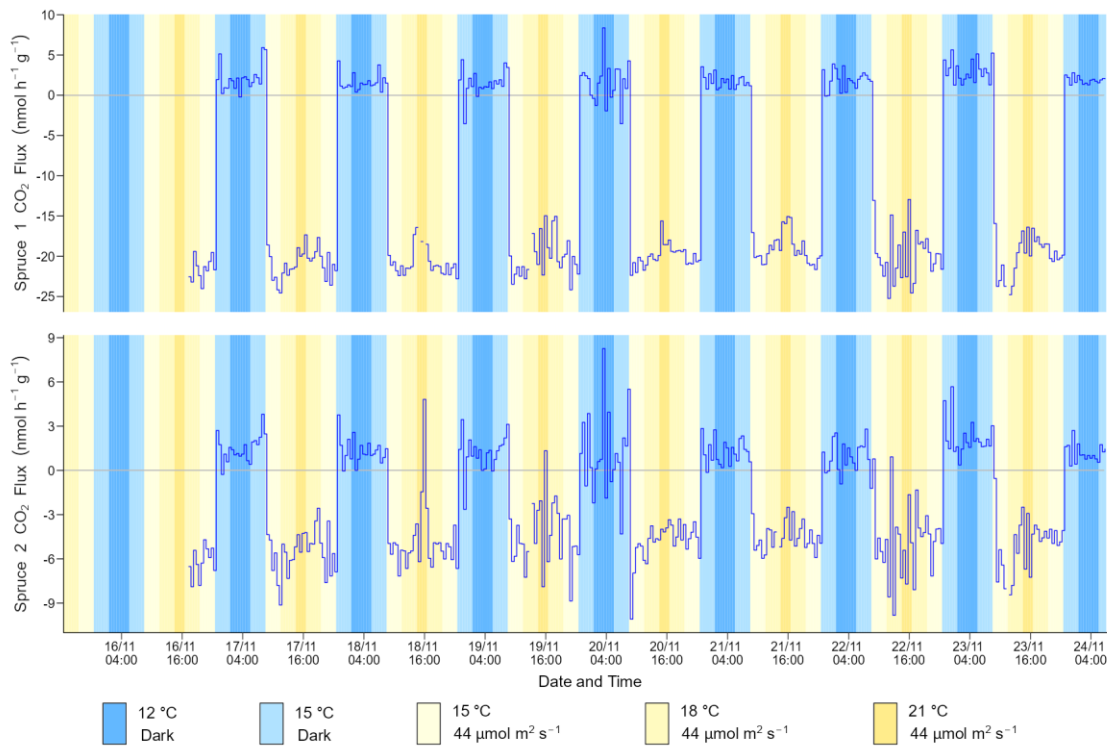


Figure S7 Time series of the CO_2 flux for Spruce 1 (top) and Spruce 2 (middle) during the Temperature Cycle.

Table S4 Chlorophyll fluorescence measurements for Spruce 1 and Spruce 2 during the Temperature Cycle. Each value is the average of three measurements taken from three branches on the same tree.

Date	Light Intensity $\mu\text{mol m}^{-2} \text{s}^{-1}$	Temperature $^{\circ}\text{C}$	Spruce 1	Spruce 2
16/11/2021	0	12	0.815	0.810
16/11/2021	44	18	0.762	0.773
19/11/2021	0	12	0.821	0.831
19/11/2021	44	18	0.820	0.785
22/11/2021	0	12	0.815	0.760
22/11/2021	44	18	0.785	0.805

The F_v/F_m values obtained for Spruce 1 and Spruce 2 during the *Temperature Cycle* are listed in Table S2.5. The values in the dark were centred around 0.8 and were similar to those observed during the *Daily Cycle*. One branch on Spruce 2 had lower measurements, and may indicate that Spruce 2 was stressed.

Light Cycle

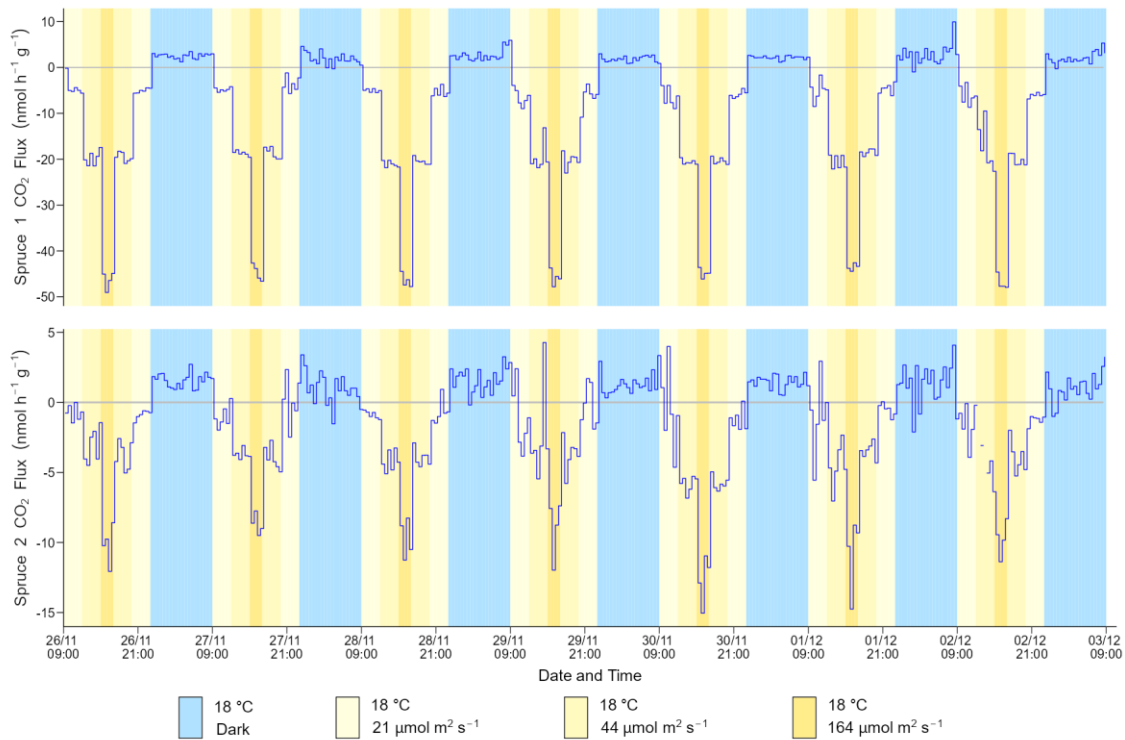


Figure S8 Time series of CO₂ flux for Spruce 1 (top) and Spruce 2 (middle) during the Light Cycle.

Table S5 Chlorophyll fluorescence measurements for Spruce 1 and Spruce 2 during the Light Cycle. Each value is the average of three measurements taken from three branches on the same tree.

Date	Light Intensity $\mu\text{mol m}^{-2} \text{s}^{-1}$	Temperature $^{\circ}\text{C}$	Spruce 1	Spruce 2
28/11/2022	0	18	0.804	0.828
28/11/2022	164	18	0.653	0.742
30/11/2021	0	18	0.694	0.768
30/11/2021	164	18	0.718	0.751
02/12/2021	0	18	0.726	0.769
02/12/2021	164	18	0.750	0.724

The F_v/F_m values recorded during the *Light Cycle* for Spruce 1 and Spruce 2 were slightly lower than those measured during the other cycles. This indicates that although Spruce 2 was stressed, it was not suffering any additional stress during the *Light Cycle*.

Standardisation

Table S6 BVOC emission fluxes from Spruce 1 standardised to 1000 $\mu\text{mol m}^{-1} \text{s}^{-1}$ and 30°C in $\mu\text{g h}^{-1}\text{s}^{-1}$. BVOC emissions ordered in order of decreasing standardised emission flux.

Formula	BVOC	Standardisation Method	Standardised Emission Flux $\mu\text{g h}^{-1} \text{s}^{-1}$
C ₁₀ H ₁₆ O	Piperitone	Combined	17.290
C ₅ H ₈	Isoprene	Biosynthesis	6.306
C ₁₀ H ₁₆	Monoterpene	Combined	0.925
C ₁₀ H ₁₄ O		Combined	0.859
C ₅ H ₈ O ₂		Biosynthesis	0.467
C ₁₀ H ₁₆ O ₂		Combined	0.455
C ₉ H ₈ O	(E)-Cinnamaldehyde	Biosynthesis	0.334
C ₁₃ H ₃₀ O ₅		Combined	0.276
C ₁₂ H ₂₀ O		Biosynthesis	0.271
C ₁₀ H ₁₆ O	Camphor	Combined	0.199
C ₅ H ₁₀ O ₆		Biosynthesis	0.196
C ₂₀ H ₂₄		Combined	0.177
C ₁₀ H ₁₄ O ₂		Combined	0.166
C ₁₀ H ₁₀		Biosynthesis	0.155
C ₁₀ H ₂₀ O ₂	Isopentyl isovalerate	Biosynthesis	0.148
C ₈ H ₁₂		Combined	0.147
C ₁₄ H ₁₆ O		Combined	0.138
C ₇ H ₁₂ N ₂		Biosynthesis	0.127
C ₃ H ₈ N ₂ O		Biosynthesis	0.082
C ₆ H ₁₄ O ₃		Combined	0.082
C ₁₂ H ₁₆		Biosynthesis	0.081
C ₁₁ H ₁₄ O		Combined	0.077
C ₂₀ H ₂₄ O ₂		Combined	0.074
C ₇ H ₄ O ₄		Combined	0.067
C ₇ H ₈ O		Combined	0.052
C ₁₇ H ₂₆ O		Combined	0.046
C ₇ H ₈		Combined	0.040
C ₉ H ₁₂		Combined	0.030
C ₁₃ H ₁₂ O		Biosynthesis	0.025
C ₁₀ H ₁₈ O ₄		Combined	0.020
C ₁₀ H ₁₄ O ₅		Combined	0.016

C ₅ H ₁₀ NO ₂		Combined	0.015
C ₁₈ H ₃₄ O		Combined	0.014
C ₃ H ₉ N		Combined	0.012

Table S7 BVOC emission fluxes from Spruce 2 standardised to 1000 $\mu\text{mol m}^{-1} \text{s}^{-1}$ and 30°C in $\mu\text{g h}^{-1}\text{s}^{-1}$. BVOC emissions ordered in order of decreasing standardised emission flux.

Formula	BVOC	Standardisation Method	Standardised Emission Flux $\mu\text{g h}^{-1} \text{s}^{-1}$
C ₁₀ H ₁₆	Monoterpene	Combined	1.226
C ₁₀ H ₁₆ O	Piperitone	Combined	1.029
C ₆ H ₈ O		Pooled	0.192
C ₂₀ H ₂₄		Pooled	0.134
C ₉ H ₈ O	(<i>E</i>)-Cinnamaldehyde	Combined	0.120
C ₆ H ₁₀ O	<i>E</i> -2-Hexenal	Pooled	0.104
C ₁₅ H ₂₄	<i>E</i> - β -Farnesene	Combined	0.097
C ₈ H ₁₂ O		Combined	0.073
C ₁₄ H ₁₆ O		Combined	0.071
C ₁₀ H ₁₄ O ₂		Pooled	0.066
C ₁₀ H ₁₆ O	Camphor	Combined	0.061
C ₇ H ₈ O		Combined	0.056
C ₁₃ H ₃₀ O ₅		Combined	0.055
C ₈ H ₁₂		Combined	0.051
C ₉ H ₁₂ O		Combined	0.046
C ₉ H ₁₂		Combined	0.035
C ₈ H ₈ O ₃	Methyl salicylate	Combined	0.033
C ₂₀ H ₂₄ O ₂		Combined	0.033
C ₇ H ₈		Combined	0.031
C ₁₀ H ₁₆		Pooled	0.029
C ₁₀ H ₁₈ O ₄		Combined	0.025
C ₆ H ₈		Combined	0.023
C ₁₀ H ₁₈ O	Eucalyptol	Combined	0.022
C ₁₇ H ₂₆ O		Combined	0.018
C ₉ H ₁₀ O ₂		Combined	0.018
C ₁₂ H ₁₆ O		Combined	0.012
C ₇ H ₁₀ O		Combined	0.009
C ₁₀ H ₁₆ O ₃		Combined	0.008

C ₈ H ₁₂		Combined	0.007
C ₆ H ₈ O		Combined	0.007
C ₁₀ H ₁₄ O ₂		Pooled	0.006
C ₁₈ H ₃₄ O		Combined	0.003

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