

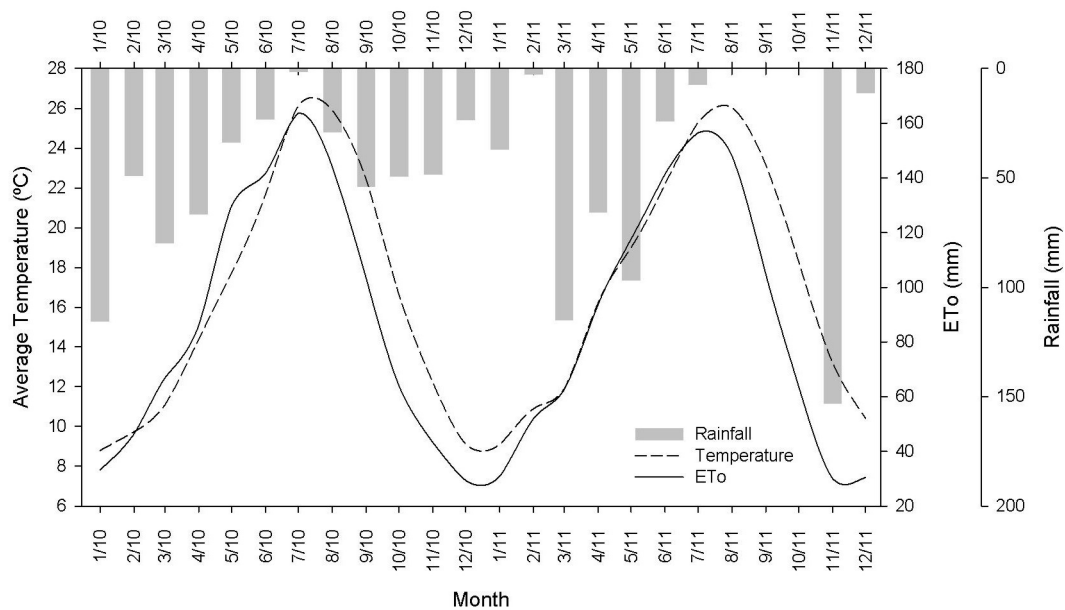
**Volatile compounds and phenolic composition of skins and seeds of 'Cabernet Sauvignon' grapes under different deficit irrigation regimes**

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Supplementary Fig. 1: Average monthly temperature (°C), total monthly reference evapotranspiration,  $ET_0$  (mm) and monthly rainfall (mm) in the 'Cabernet Sauvignon' vineyard during the experimental period (2010-2011).

Supplementary Table 1

Post-veraison water stress integral (MPa\*day) in the rainfed application and in the treatments watered with 25, 50 and 75 % of the estimated crop evapotranspiration ( $ET_c$ ) during the post-veraison period

Year	T			
	Rainfed	0.25 $ET_c$	0.50 $ET_c$	0.75 $ET_c$
2010	35.00	28.33	22.85	22.88
2011	30.02	24.54	19.52	18.15

## Supplementary Table 2

Volatile compounds identified in grape samples during 2010 and 2011

Volatile compound	RT	RI	ID
Alcohols			
2-Butanol	9.33-10.24	1039-1067	MS;RI;B;St
1-Propanol	9.84-10.81	1055-1085	MS;RI;B
2-Methyl-1-propanol	11.85-12.99	925-1127	MS;RI;B;St
1-Butanol	13.87-15.17	950-1182	MS;RI;B
3-Methyl-1-butanol	16.14-17.54	1045-1120	MS;RI;B
1-Pentanol	17.78-19.25	1061-1240	MS;RI;B;S
1-Hexanol	21.60-23.14	1087-1286	MS;RI;B;S
(Z) 3-Hexen-1-ol	22.9-24.44	1107-1358	MS;RI;B;S
(E) 2-Hexen-1-ol	23.17-25.21	1121-1425	MS;RI;B
2-Octanol	23.88-25.42	1413-1458	MS;RI;B
1-Heptanol	25.29-26.84	1453-1498	MS;RI;B;S
2-Ethyl-1-hexanol	26.48-28.04	1487-1534	MS;RI;B
Aldehydes			
2-Methyl-propanal	4.04-4.26	814-829	MS
2-Methyl-butanal	5.97-6.39	826-922	MS;RI
3-Methyl-butanal	6.07-6.50	833-927	MS;RI;B;S
Hexanal	11.32-12.40	921-1112	MS;RI;B;S
Heptanal	15.22-16.54	1201-1217	MS;RI;B;S
(E) 2-Hexenal	16.97-18.32	1052-1229	MS;RI;B;St
Esters			
Methyl acetate	4.30-4.55	831-846	MS
Ethyl propionate	7.10-7.70	961-986	MS;RI;B
Ethyl 2-methylpropionate	7.25-7.89	968-993	MS;RI;B
Ethyl butanoate	9.56-10.48	1046-1075	MS;RI;B
Ethyl 3-methylbutanoate	10.66-11.68	1080-110	MS;RI
Hexyl acetate	18.45-19.91	1240-1258	MS;RI;B
Ethyl octanoate	24.37-25.89	1427-1472	MS;RI;B
Terpenes			
D-limonene	15.19-16.49	1201-1217	MS;RI;B;S
3,7-Dimethyl- 1,6-octadien-3-ol	29.06-30.65	1565-1615	MS;RI;B;St

RT (retention time); RI (Linear retention indices calculated). ID: Method of identification (MS-comparison with mass spectrum stored in NIST library; RI-comparison of linear retention indices with the literature; B- comparison of other authors; St-comparison of retention time and mass spectrum with those of authentic standards).