

Supporting Information

Phenylalkanoid Glycosides (Non-Salicinoids) from Wood Chips of *Salix triandra* × *dasyclados* Hybrid Willow

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Table S2. NMR chemical shift data of chavicol- α -L-arabinofuranosyl-(1'' \rightarrow 6')- β -D-glucopyranoside (**18**) and chavicol-- β -D-apiofuranosyl-(1'' \rightarrow 6')- β -D-glucopyranoside (**20**)

Table 1. NMR chemical shift data of *p*-hydroxybenzylacetone- β -D-glucopyranoside (**5**).

Pos ⁿ	δ_c^a	δ_H^a	J_{H-H} (Hz); <i>multiplicity</i> ^b	COSY	HMBC	$\delta_c^{c,d}$	$\delta_H^{c,d}$	J_{H-H} (Hz); <i>multiplicity</i> ^{b,c,d}
1	158.0	-	-	-	-	157.1	-	-
2	119.7	7.07	8.6; <i>d</i>	H-3	C-1, C-3/5, C-4	118.0	7.11	8.4; <i>d</i>
3	132.5	7.23	8.6; <i>d</i>	H-2	C-1, C-2/6, C-7	130.1	7.00	8.4; <i>d</i>
4	138.6	-	-	-	-	136.2	-	-
5	132.5	7.23	8.6; <i>d</i>	H-6	C-1, C-2/6, C-7	130.1	7.00	8.4; <i>d</i>
6	119.7	7.07	8.6; <i>d</i>	H-5	C-1, C-3/5, C-4	118.0	7.11	8.4; <i>d</i>
7	31.4	2.86	<i>m</i>	-	C-3/5, C-4, C-8, C-9	30.2	2.78	<i>m</i>
8	47.5	2.88	<i>m</i>	-	C-4, C-7, C-9	46.2	2.78	<i>m</i>
9	218.5	-	-	-	-	211.0	-	-
10	32.3	2.19	<i>s</i>	-	C-9	30.2	2.11	<i>s</i>
1'	103.4	5.05	7.7; <i>d</i>	H-2'	C-1	102.4	4.87	7.8; <i>d</i>
2'	75.9	3.53	7.8; <i>d</i>	H-1'	C-3'	75.1	3.39-4.40	<i>m</i>
3'	78.6	3.58	9.0; <i>d</i>	-	C-2', C-4'	78.3	3.39-4.40	<i>m</i>
4'	72.6	3.47	9.0	-	C-5'	71.6	3.39-4.40	<i>m</i>
5'	78.9	3.43	<i>m</i>	H-6'	C-4'	78.2	3.39-4.40	<i>m</i>
6'	63.7	3.92	12.4, 2.1; <i>dd</i>	H-5'	C-5'	62.7	3.39-4.40	<i>m</i>
		3.73	12.4, 5.8; <i>dd</i>					

^a Data collected in 80:20 D₂O:CD₃OD (4:1). Spectra were referenced to TSP-d4 at δ 0.00.

^b *s* singlet; *d* doublet; *dd* double doublet; *m* multiplet.

^c Data collected in CD₃OD.

^d Gao, L.; Xu, X.; Yang, J. Chemical constituents of the roots of *Rheum officinale*. *Chem. Nat. Compd.* **2013**, *49*, 603–605, doi:10.1007/s10600-013-0689-7.

Table S2: NMR chemical shift data of chavicol- α -L-arabinofuranosyl-(1'' \rightarrow 6')- β -D-glucopyranoside (**18**) and chavicol-- β -D-apiofuranosyl-(1'' \rightarrow 6')- β -D-glucopyranoside (**20**)

Pos ⁿ	18					20				
	δ_c	δ_H	J_{H-H} (Hz); <i>multiplicity</i> ^a	COSY	HMBC	δ_c	δ_H	J_{H-H} (Hz); <i>multiplicity</i> ^a	COSY	HMBC
1	158.1	-	-	-	-	158.1	-	-	-	-
2	119.9	7.07	8.7; <i>d</i>	H-3/5	C-1, C-4, C-3/5	119.8	7.07	8.6; <i>d</i>	H-3/5	C-1, C-4, C-3/5
3	132.9	7.22	8.7; <i>d</i>	H-2/6	C-1, C-2/6	132.8	7.22	8.6; <i>d</i>	H-2/6	C-1, C-2/6, C-7
4	138.5					138.4				
5	132.9	7.22	8.7; <i>d</i>	H-2/6	C-1, C-2/6	132.8	7.22	8.6; <i>d</i>	H-2/6	C-1, C-2/6, C-7
6	119.9	7.07	8.7; <i>d</i>	H-3/5	C-1, C-4, C-3/5	119.8	7.07	8.6; <i>d</i>	H-3/5	C-1, C-4, C-3/5
7	41.7	3.36	6.7; <i>d</i>	H-8		41.6	3.36	6.9; <i>d</i>	H-8	C-2/6, C-3/5, C-4, C-8
8	141.5	6.01	16.9, 10.1, 6.7; <i>ddt</i>	H-7, H-9		141.3	6.01	16.9, 10.1, 6.7; <i>ddt</i>	H-7, H-9	C-4, C-7
9	118.6	5.07	<i>m</i>	H-8		118.4	5.08	<i>m</i>	H-8	C7, C8
1'	103.5	5.04	7.7; <i>d</i>	H-2'	C-1, C-4'	103.5	5.03	7.6; <i>d</i>	H-2'	C-1, C-2'
2'	76.0	3.54	7.7; <i>d</i>	H-1'	C-1'	78.5	3.53	7.6; <i>d</i>	H-1', H-3'	C-1', C-3'
3'	72.7	3.50	9.1; <i>d</i>		C-1'	76.1	3.56	8.9; <i>d</i>	H-2'	C-4'
4'	78.6	3.57	9.0; <i>d</i>		C-3'	72.6	3.48	9.2; <i>d</i>		C-5', C-6'
5'	78.1	3.73	<i>m</i>			77.9	3.7	<i>m</i>	H-6'	
6'	69.8	4.04	11.1; <i>d</i>		C-3', C-1''	70.7	4.02	2.3, 11.4; 3.72 <i>dd</i> <i>m</i>	H-5'	C-4', C-5', C-1''
1''	111.2	5.00	1.3; <i>d</i>	H-2''	C-6', C-3'', C-4''	112.2	5.05	3.1; <i>d</i>	H-2''	C-6', C-2'', C-3'', C-4''
2''	84.2	4.06	1.5, 3.3; <i>dd</i>	H-1'', H-3''	C-1'', C-3''	79.6	3.94	3.1; <i>d</i>	H-1''	C-1'', C-5''
3''	79.7	3.89	3.3, 5.9; <i>dd</i>	H-2''	C-2'', C-4'', C-5''	82.3	-	-	-	-
4''	87.0	4.00	3.3; 5.9; <i>dd</i>		C-2'', C-3'', C-5''	76.6	3.83	10.1; <i>d</i> 3.99 10.1; <i>d</i>	H-4'', H-5''	C-1'', C-2'', C-3'', C-5''
5''	64.1	3.74	3.3, 12.3; <i>dd</i> 3.64 5.7, 12.3; <i>dd</i>		C-3'', C-4''	66.5	3.6	<i>s</i>	H-4''	C-2'', C-3'', C-4''

Data collected in 80:20 D₂O:CD₃OD (4:1). Spectra were referenced to TSP-d4 at δ 0.00.

^a *s* singlet; *d* doublet; *dd* double doublet; *m* multiplet.