

DOI: <http://dx.doi.org/10.21577/1984-6835.20230050>

***In Vitro* Antimollicute Properties of *Rosmarinus officinalis* Extract**

Propriedades Antimollicute *In Vitro* do Extrato de *Rosmarinus officinalis*

Debora H. Mafra,^a (<https://orcid.org/0000-0002-9721-6398>) Italo R. Garcia,^b

(<https://orcid.org/0000-0002-8934-690X>) Giovana Vieira,^c (<https://orcid.org/0000-0001-9692-8070>) Alessandro Guedes,^d (<https://orcid.org/0000-0001-7785-8281>) Caio

M. M. De Cordova^{e,*} (<https://orcid.org/0000-0001-6090-0367>)

^a Universidade de Blumenau, Campus 1, Programa de pós-graduação em Química, Rua Antônio da Veiga 140, CEP 89030-903, Blumenau-SC, Brasil.

^b Universidade de Blumenau, Campus 3, Curso de Biomedicina, Rua São Paulo 2171, lab. A-503, CEP 89030-001, Blumenau-SC, Brasil.

^c Universidade de Blumenau, Campus 1, Departamento de Química, Rua Antônio da Veiga 140, CEP 89030-903, Blumenau-SC, Brasil.

^d Universidade de Blumenau, Campus 3, Departamento de Ciências Farmacêuticas, Rua São Paulo 2171, CEP 89030-001, Blumenau-SC, Brasil.

^e Universidade de Blumenau, Campus 3, Rua São Paulo 2171, Departamento de Ciências Farmacêuticas, lab. A-503, CEP 89030-001, Blumenau-SC, Brasil.

[*cmcordova@furb.br](mailto:cmcordova@furb.br)

Supplementary Information

Table 6S. ^{13}C NMR data of the compound Ro1 and comparison with literature values . Note: There is a noticeable difference in the value of C1 (CH₂) of Ro1 with the reported value, however the displacement value of Ro1 is compatible with the value of C1 of betulinic acid.

C position	C type	Ro1 δC	Betulinic acid δC (Shai et al, 2008)
1	CH ₂	38.9929	31.9
2	CH ₂	28.5727	28.1
3	CH	77.8043	78.1
4	C	39.2909	39.4
5	CH	55.9040	55.7
6	CH ₂	19.9953	18.6
7	CH ₂	34.9368	34.7
8	C	ND	40.9
9	CH	50.9667	50.8
10	C	37.7618	37.3
11	CH ₂	21.6219	21.1
12	CH ₂	26.0989	25.9
13	CH	38,6430	38.4
14	C	44.6947	42.4
15	CH ₂	31.1269	31.1
16	CH ₂	32.7338	32.7
17	C	56.4612	56.3
18	CH	47.6623	47.6
19	CH	49.5672	49.5
20	C	151.3840	150.7
21	CH ₂	30.2327	30.1
22	CH ₂	37.3730	37.5
23	CH ₃	29.1313	28.5
24	CH ₃	16.9620	16.3
25	CH ₃	16.8469	16.3
26	CH ₃	16.7625	16.2

27	CH3	15.4209	14.8
28	C	178.2993	178.7
29	CH2	110.6936	110.3
30	CH3	18.9846	19.4

ND: The C-8 signal was was sidden by the solvent signal. δ : Chemical displacement.

L.J. Shai, L.J. McGaw, M.A. Aderogba, L.K. Mdee, J.N. Eloff.

Four pentacyclic triterpenoids with antifungal and antibacterial activity from *Curtisia dentata* (Burm.f) C.A. Sm. Leaves, Journal of Ethnopharmacology 2008;119:238-44.

<https://doi.org/10.1016/j.jep.2008.06.036>

Figure 1S. Infrared (a), ^{13}C NMR (b), ^1H NMR (c) and ^{13}C DEPT (d) spectra of the compound Ro1.



