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MOLECULAR NETWORKING-BASED APPROACH OF ALOE DJIBOUTIENSIS, ANTIOXIDANT ACTIVITY AND IN VIVO TOXICITY OF THIS ENDEMIC SPECIES IN DJIBOUTI



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INTRODUCTION

Aloe djiboutiensis is a native plant in Djibouti, named botanically in 2007. The leaves and the latex are used traditionally used in Djibouti. The present research work was conducted to assess the possible toxic effect and antioxidant the latex and the leave extract, as well as to detect rapidly the previously characterized known molecules in the genus Aloe and to create a metabolic fingerprinting of A. djiboutiensis [1].

RESULTS AND DISCUSSION

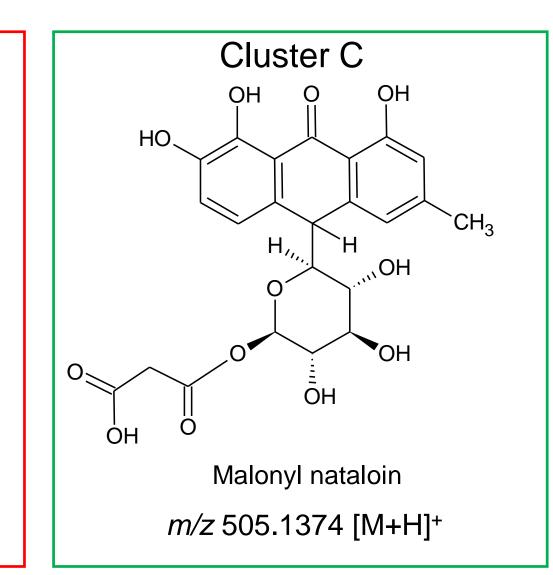
The results for antioxidant activity, the *in vitro* and *in vivo* toxicity are reported in the figure. The latex (ADL) has the high level of phenol content (83.01 \pm 0.8 mg GAE/g DW) and moderate level of flavonoid content (13.12 \pm 0.11 mg QE/g DW).

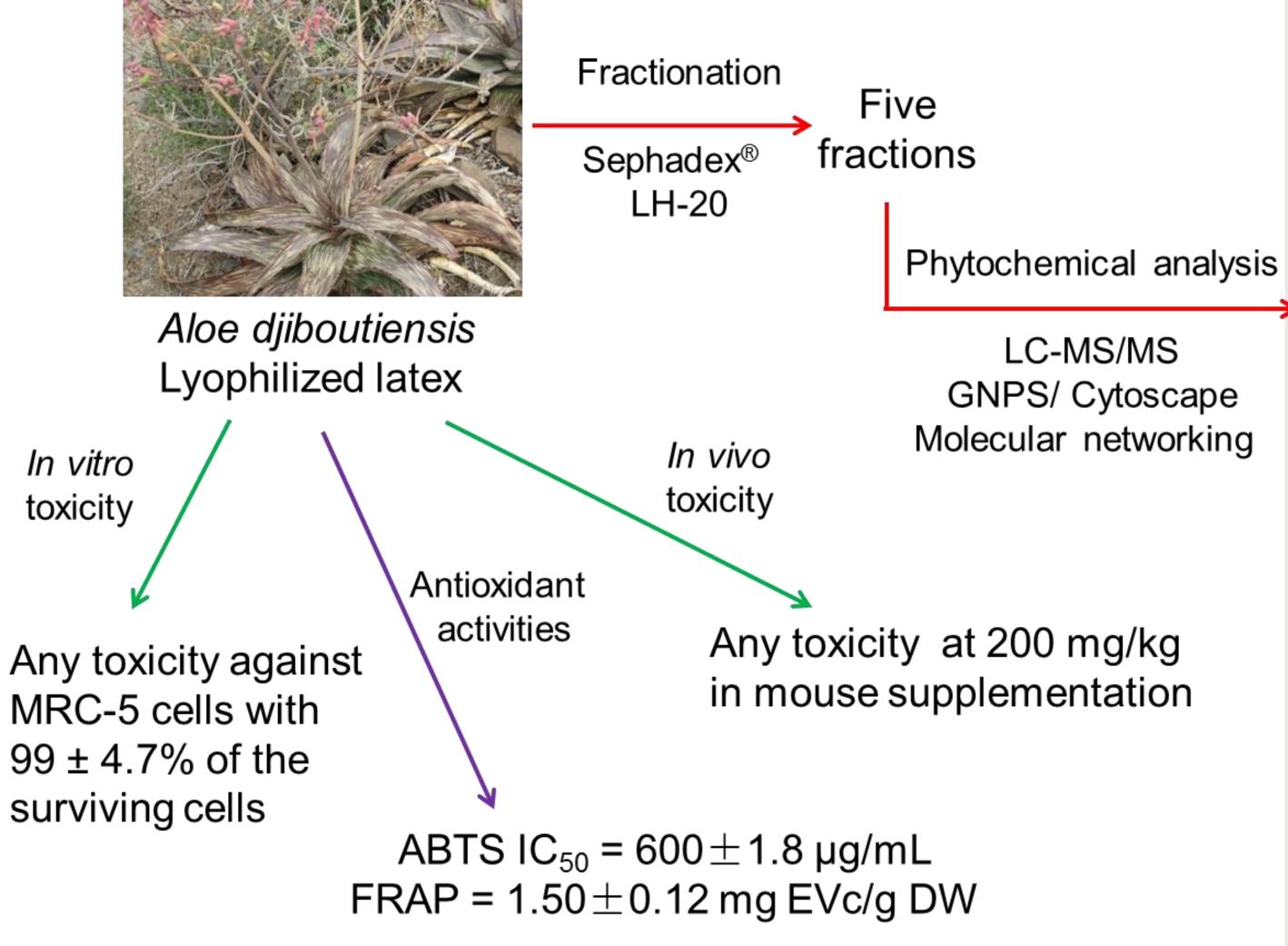
The latex and five fractions of *Aloe djiboutiensis*, obtained after LH-20 fractionation of latex, were analyzed, and the phytochemical profile is evaluated using LC-MS/MS analysis. The identification/annotation of the chemical compounds is based on the comparison of commercial standards, the chemicals present in our in-house database (DataNat database, n° IDDN: FR.001.480019.000.S.P.2020.000.10300) and

Global Natural Products Social Molecular Networking (GNPS) [2].

GNPS is used for the generation of molecular networking, and Cytoscape software 3.8.0 is used for the visualization of them. In total, 712 nodes are visible in molecular networking, which included 67 clusters (node>2). The observation of molecular network reveals the presence of eleven (11) annotated clusters named, respectively A, B, C, D, E, F, G, H, I, L, and M.

Cluster B Cluster A Aloinoside A/B Aloin A/B *m*/*z* 565.1915 [M+H]⁺ m/z 419.1441 [M+H]+



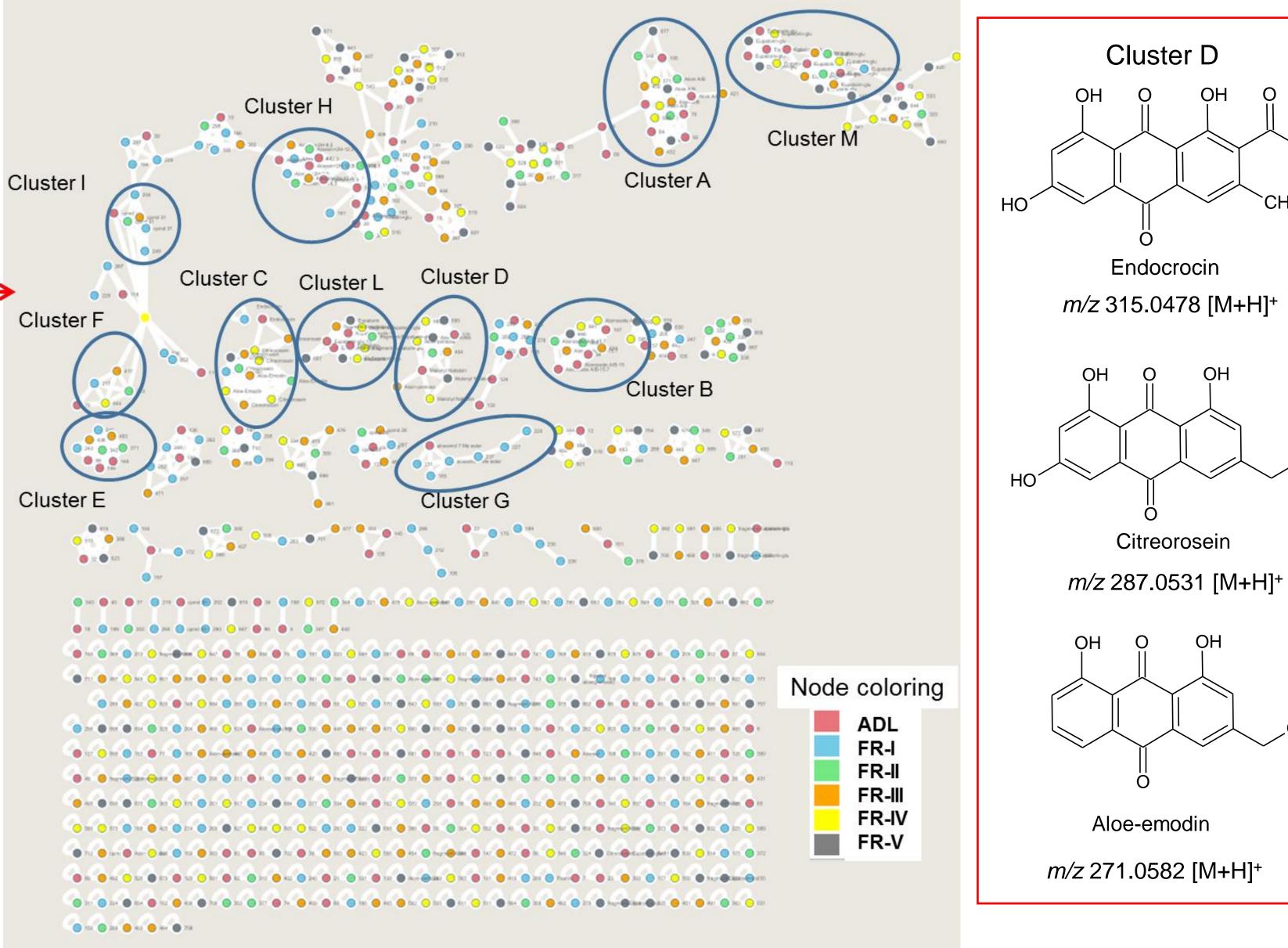


CONCLUSION

According to the obtained results, the traditional use of the latex of A. djiboutiensis is safe until at less 18 mg/kg for human level and has beneficial effects for human health (following equations for dose conversion between animals and human).

In the figure are visible the relevant structures annotated in the molecular network.

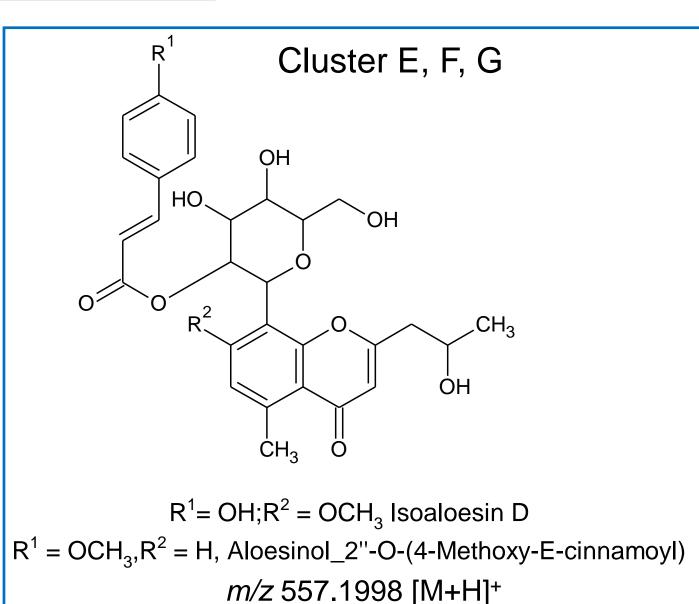
In total twenty one (21) compounds are identified: nine (9) anthraquinones derivatives, ten (10) chromones derivatives, two (2) flavonols/ chromones isomers.



Cluster L

CH₃

$$H_3C$$
 CH_3
 CH



REFERENCES

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2) Mingxun Wang, Jeremy J. Carver, Vanessa V. Phelan, Laura M. Sanchez, Neha Garg, Yao Peng, Don Duy Nguyen et al.; Nature biotechnology 34, no. 8 (2016): 828. PMID: 27504778