

Supporting Information

Figure Legends

Figure S1.

FT infra red spectrum of crude acids showing two absorptions attributed to carbonyl stretches of acids (1708 cm^{-1}) and esters (1726 cm^{-1})

Figure S2.

Electron ionization mass spectrum of the putative C_{80} hexacyclic triacid mono'octyl'ester ('the unknown').

Figure S3.

a. Positive ion Orbitrap electrospray mass spectrum of the trimethyl ester of the putative C_{80} hexacyclic triacid mono'octyl'ester ('the unknown'). A formula of $\text{C}_{91}\text{H}_{164}\text{O}_8\text{NH}_4$ is supported by the mass of the ammoniated molecular ion (1403.2749 observed; 1403.2764 required).

b. Negative ion Orbitrap electrospray mass spectrum of the putative C_{80} hexacyclic triacid mono'octyl'ester ('the unknown'). A formula of $\text{C}_{88}\text{H}_{157}\text{O}_8$ is supported by the mass of the deprotonated molecular ion (1342.1857 observed; 1342.1884 required).

Figure S4.

a. Electron ionization mass spectrum and GC retention time of the alcohol (TMS ether) obtained by saponification of the putative C_{80} hexacyclic triacid mono'octyl'ester ('the unknown'), followed by examination by GC-MS.

b. Electron ionization mass spectrum and GC retention time of authentic 2-ethylhexanol (TMS ether).

c. Electron ionization mass spectrum and GC retention time of *n*-octanol (TMS ether).

Figure S1.

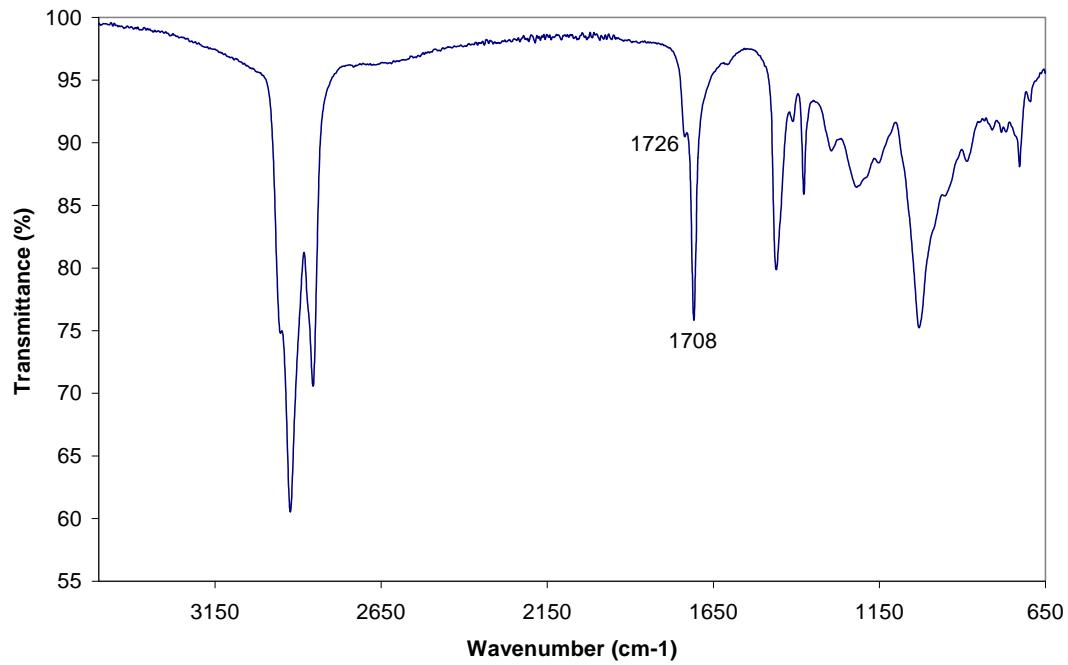


Figure S2.

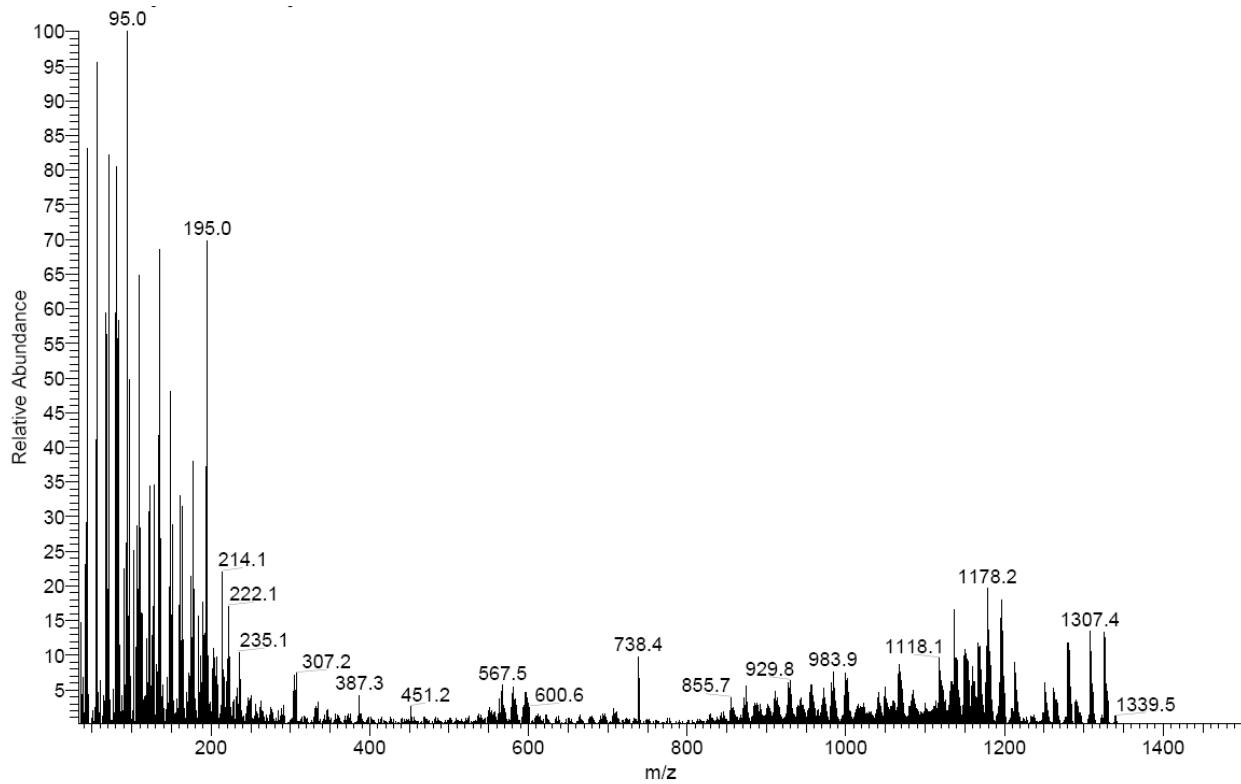


Figure S3.

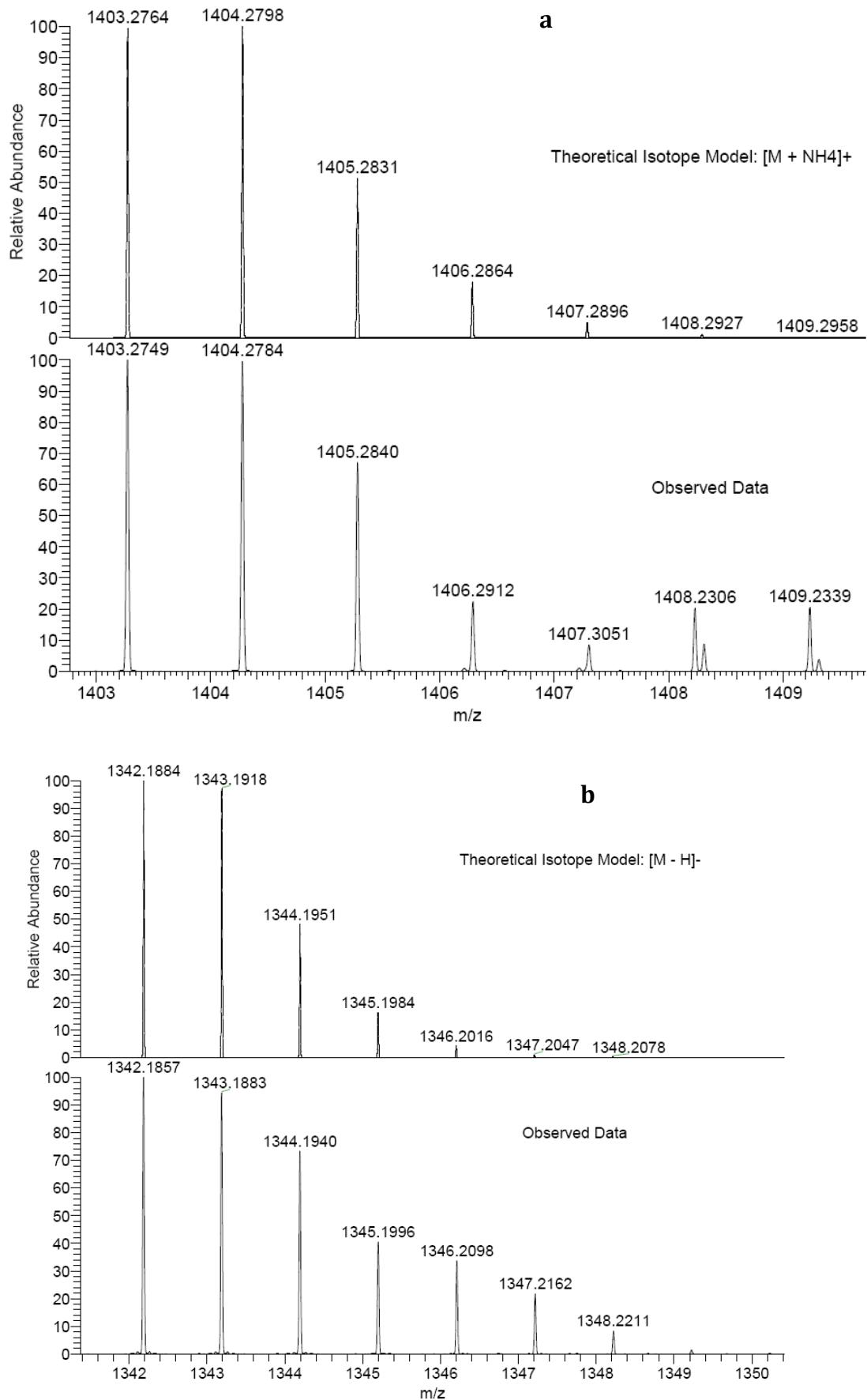
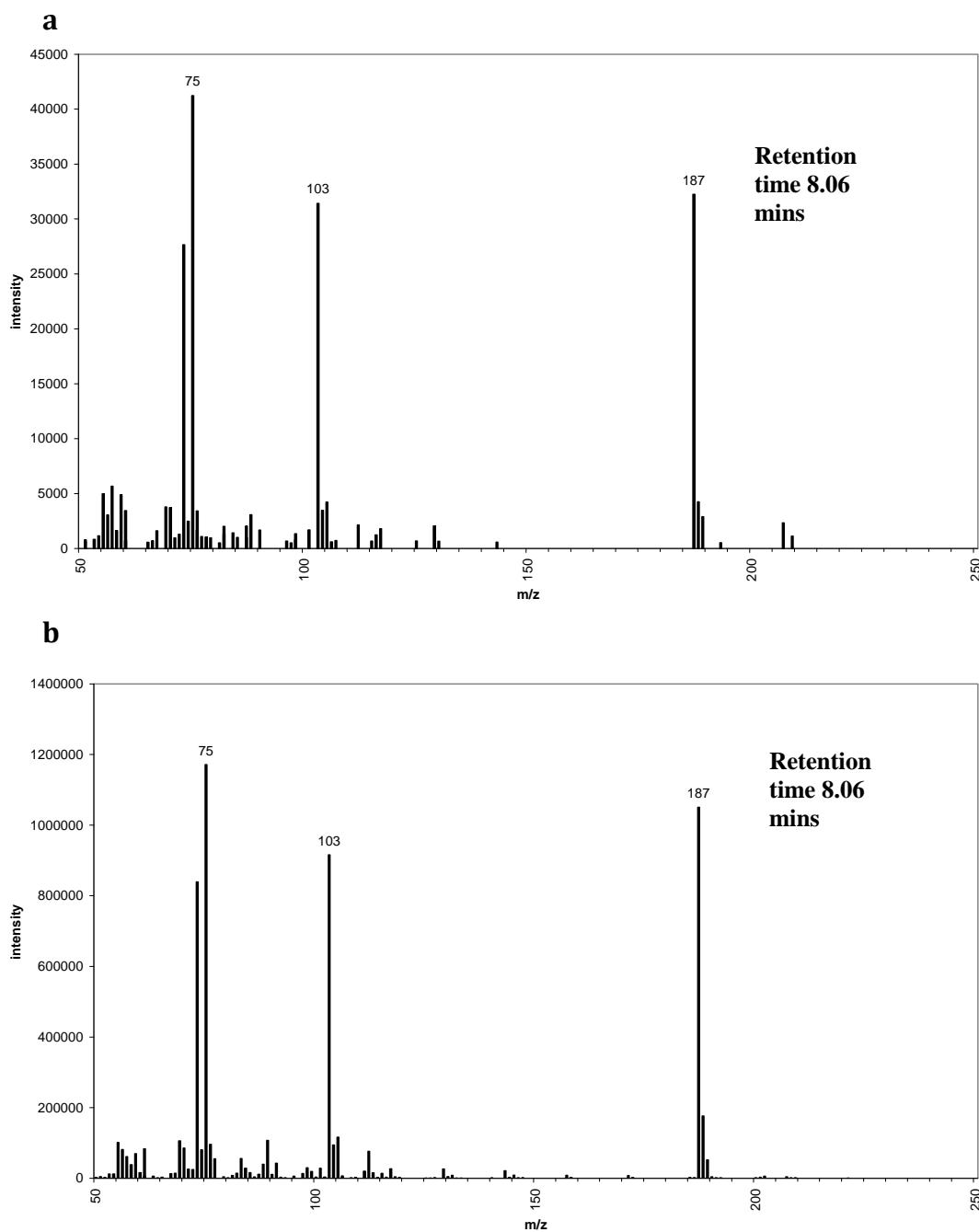


Figure S4.



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