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ALKALOIDS OF THE MONGOLIAN FLORA

I. ALKALOIDS OF THE EPIGEAL PART OF *Aconitum barbatum*

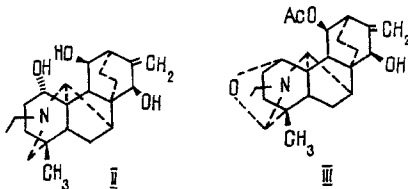
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Continuing the separation of the total alkaloids of the epigeal part of *Aconitum barbatum* Pers. [1, 2], we have isolated three bases.

Alkaloid (I), with the composition $C_{25}H_{41}NO_7$, was identified on the basis of a study of its spectral characteristics and comparison with an authentic specimen as delsoine [3].

Alkaloid (II) had the composition $C_{22}H_{33}NO_3$, mp 191-193°C (acetone). On acetylation it gave a triacetyl derivative. A study of the spectral characteristics of the alkaloid and of its triacetate enabled us to suggest a structure [4] coinciding with that of the alkaloid lepenine [5]. A direct comparison of the IR spectra of lepenine and of the alkaloid that had been isolated showed their identity. A study of the ^{13}C NMR spectrum of lepenine, which was not taken into consideration in the proof of its structure [5], agreed with this. The multiplicities of the signals were determined from the "off-resonance" spectrum, and the assignment of the signals was made by comparison with the spectrum of 11-acetyl-1,19-epoxydenudatine (III) [2].



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TABLE 1. Details of the ^{13}C NMR Spectra (100 MHz) of Lepenine (II) ($\text{C}_5\text{D}_5\text{N}$) and of 11-Acetyl-1,19-Epoxydenudatine (III) (CDCl_3)

Carbon	II	III	Carbon	II	III
1	70.1	68.3	12	42.4	43.3
2	31.7	24.1	13	23.8	24.3
3	39.3	29.7	14	25.1	26.9
4	33.8	37.5	15	77.9	77.1
5	52.8	49.6	16	155.5	153.7
6	28.2	24.4	17	108.7	110.7
7	47.9	47.4	18	23.3	18.6
8	44.3	45.5	19	51.1	92.9
9	54.3	46.5	20	68.3	69.8
			N		
10	51.6	49.4	CH ₂	57.3	48.4
11	73.2	74.2	CH ₃	13.8	14.1

Alkaloid (III) had the composition $\text{C}_{24}\text{H}_{35}\text{NO}_4$, mp 137-139°C, and differed from lepenine by the presence of an acetoxy group in place of a hydroxy group. Alkaline hydrolysis gave an amino alcohol identical with lepenine. The presence in PMR spectrum of the signal of C-11- α proton at 5.52 ppm (d, $J = 9$ Hz) that was shifted downfield in comparison with the analogous signals in the spectrum of lepenine showed that the acetoxy group was located at C-11. Consequently, this alkaloid was lepetine [5].

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