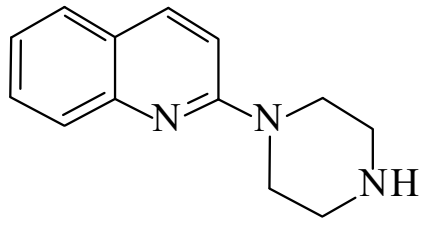
>>**Entry Name: 2-(1-Piperazinyl)quinoline,** 9CI, 8CI
Synonym(s): Quipazine, INN



CAS Registry Number: 4774-24-7

Molecular Formula: $C_{13}H_{15}N_3$
Molecular Weight: 213.282
Accurate Mass: 213.126597
Percentage Composition: C 73.21%; H 7.09%; N 19.70%
Use/Importance: Non-selective 5-HT-receptor agonist. Antiparkinsonian drug, antidepressant, oxytocic agent. Pharmacological tool
Melting Point: Mp 82-83°
Calculated Log P Data: Log P 2.25 (calc)

>>**Derivative:** Maleate (1:1)
Synonym(s): Quipazine maleate, USAN**.** MR 1291

CAS Registry Number: 5786-68-5
Hazard & Toxicity: LD_{50} (mus, orl) 225 mg/kg

References:

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