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| Name | ANDARINE |
| Structure | |
| Systematic name | (2S)-3-(4-Acetamidophenoxy)-2-hydroxy-2-methyl-N-[4-nitro-3-(trifluoromethyl)phenyl]propanamide |
| Formula | C ₁₉ H ₁₈ F ₃ N ₃ O ₆ |
| MW | 441.3579 |
| Monoisotopic mass | 441.1148 |
| Mp | 70-74°C |
| H bond acceptors | 9 |
| H bond donors | 3 |
| Acid pKa | 12 (p-nitroanilide), 14 (tertiary OH) |
| Basic pka | -- |
| ACD Log D pH 5.5 | 4.01 |
| ACD Log D pH 7.4 | 4.01 |
| Solubility | Ethanol, DMSO. 1.2 mg/mL in water |
| LD50 | -- |
| Therapeutic cat | antiBPH (benign prostatic hypertrophy) |
| ATC | Investigational new drug |
| Receptors | AR partial agonist |
| Nomi commerciali (IT) | |
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