

Title (en)  
SUBSTITUTED TETRAHYDROISOQUINOLINE DERIVATIVES AS MODULATORS OF DOPAMINE D<sub>3</sub> RECEPTORS

Title (de)  
SUBSTITUIERTE TETRAHYDROISOCHINOLINEDERIVATE ALS DOPAMINE D<sub>3</sub> REZEPTOR-MODULATOREN

Title (fr)  
DERIVES DE TETRAHYDRO-ISOQUINOLINE SUBSTITUEE, UTILES EN TANT QUE MODULATEURS DES RECEPTEURS D<sub>3</sub> DE LA DOPAMINE

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Application  
**EP 98929278 A 19980427**

Priority  
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• GB 9708694 A 19970430

Abstract (en)  
[origin: WO9849145A1] Compounds of formula (I), wherein R<sub>1</sub>> represents a substituent selected from a hydrogen or halogen atom; a hydroxy, cyano, nitro, trifluoromethyl, trifluoromethoxy, trifluoromethanesulfonyloxy, pentafluoroethyl, C<sub>1</sub>-4alkyl, C<sub>1</sub>-4alkoxy, arylC<sub>1</sub>-4alkoxy, C<sub>1</sub>-4alkylthio, C<sub>1</sub>-4alkoxyC<sub>1</sub>-4alkyl, C<sub>3</sub>-6cycloalkylC<sub>1</sub>-4alkoxy, C<sub>1</sub>-4alkanoyl, C<sub>1</sub>-4alkoxycarbonyl, C<sub>1</sub>-4alkylsulphonyl, C<sub>1</sub>-4alkylsulphonyloxy, C<sub>1</sub>-4alkylsulphonylC<sub>1</sub>-4alkyl, arylsulphonyl, arylsulphonyloxy, arylsulphonylC<sub>1</sub>-4alkyl, C<sub>1</sub>-4alkylsulphonamido, C<sub>1</sub>-4alkylamido, C<sub>1</sub>-4alkylsulphonamidoC<sub>1</sub>-4alkyl, C<sub>1</sub>-4alkylamidoC<sub>1</sub>-4alkyl, arylsulphonamido, arylcarboxamido, arylsulphonamidoC<sub>1</sub>-4alkyl, arylcarboxamidoC<sub>1</sub>-4alkyl, aroyl, aroylC<sub>1</sub>-4alkyl, or arylC<sub>1</sub>-4alkanoyl group; a group R<sub>3</sub>>OCO(CH<sub>2</sub>)<sub>p</sub>, R<sub>3</sub>>CON(R<sub>4</sub>>)(CH<sub>2</sub>)<sub>p</sub>, R<sub>3</sub>>R<sub>4</sub>>NCO(CH<sub>2</sub>)<sub>p</sub> or R<sub>3</sub>>R<sub>4</sub>>NSO<sub>2</sub>(CH<sub>2</sub>)<sub>p</sub> where each of R<sub>3</sub>> and R<sub>4</sub>> independently represents a hydrogen atom or a C<sub>1</sub>-4alkyl group or R<sub>3</sub>>R<sub>4</sub>> forms part of a C<sub>3</sub>-6 azacycloalkane or C<sub>3</sub>-6(2-oxo)azacycloalkane ring and p represents zero or an integer from 1 to 4; or a group Ar<sub>1</sub>>Z, wherein Ar<sub>1</sub>> represents an optionally substituted phenyl ring or an optionally substituted 5- or 6-membered aromatic heterocyclic ring and z represents a bond, O, S, or CH<sub>2</sub>; R<sub>2</sub>> represents a hydrogen atom or a C<sub>1</sub>-4alkyl group; q is 1 or 2; Ar represents an optionally substituted phenyl ring or an optionally substituted 5- or 6-membered aromatic heterocyclic ring; or an optionally substituted bicyclic ring system; and salts thereof. Compounds of formula (I) and their salts have affinity for dopamine receptors, in particular the D<sub>3</sub> receptor, and thus potential in the treatment of conditions wherein modulation of the D<sub>3</sub> receptor is beneficial, e.g. as antipsychotic agents.

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IPC 8 full level  
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