

**United States Patent** [19]

Müller et al.

[11] **Patent Number:** 5,684,014[45] **Date of Patent:** Nov. 4, 1997[54] **CYCLOALKANO-INDOLE AND  
-AZAINDOLE DERIVATIVES**

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C07D 471/04

[52] **U.S. Cl.** ..... 514/292; 546/86; 546/87

[58] **Field of Search** ..... 546/86, 87; 514/292

[56] **References Cited**

## U.S. PATENT DOCUMENTS

3,632,807 1/1972 Maurer et al. .... 546/345  
4,775,680 10/1988 Gillard et al. .... 514/411

## FOREIGN PATENT DOCUMENTS

0 234 708 A1 9/1987 European Pat. Off. .  
0 300 676 A2 1/1989 European Pat. Off. .  
0 310 179 A2 4/1989 European Pat. Off. .  
0 496 237 A2 7/1992 European Pat. Off. .  
0509359 10/1992 European Pat. Off. .  
0 617 035 A1 9/1994 European Pat. Off. .

## OTHER PUBLICATIONS

Heterocycles, vol. 22, No. 10, 1984 (pp. 2277-2279).

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[57] **ABSTRACT**

Cycloalkano-indole and -azaindole derivatives are prepared by reaction of appropriately substituted carboxylic acids with amines. The cycloalkano-indole and -azaindole derivatives are suitable as active compounds for medicaments, preferably antiatherosclerotic medicaments.

**11 Claims, No Drawings**

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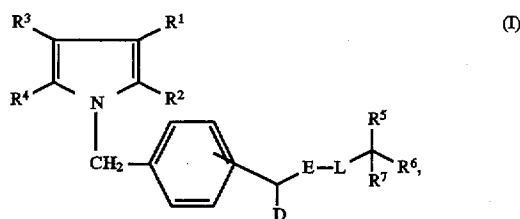
**CYCLOALKANO-INDOLE AND  
-AZAINDOLE DERIVATIVES**

The present invention relates to cycloalkano-indole and -azaindole derivatives, processes for their preparation and their use as medicaments, in particular as antiatherosclerotic medicaments.

It is known that increased blood levels of triglycerides (hypertriglyceridaemia) and cholesterol (hypercholesterolaemia) are associated with the genesis of atherosclerotic vessel wall changes and coronary heart diseases.

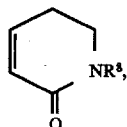
A distinctly increased risk of the development of coronary heart disease is moreover present if these two risk factors occur in combination, which is accompanied, in turn, with an overproduction of apolipoprotein B-100. There is therefore, as before, a great need to make available effective medicaments for the control of atherosclerosis and coronary heart diseases.

The present invention relates to cycloalkano-indole and -azaindole derivatives of the general formula (I)



in which

$R^1$  and  $R^2$ , including the double bond connecting them, together form a phenyl or pyridyl ring or a ring of the formula



wherein

$R^8$  denotes hydrogen or straight-chain or branched alkyl having up to 4 carbon atoms,

$R^3$  and  $R^4$ , including the double bond connecting them, together form a phenyl ring or a 4- to 8-membered cycloalkene or oxocycloalkene radical,

all ring systems mentioned under  $R^1/R^2$  and  $R^3/R^4$  optionally being substituted up to 3 times by identical or different halogen, trifluoromethyl, carboxyl or hydroxyl substituents, by straight-chain or branched alkoxy or alkoxycarbonyl each having up to 6 carbon atoms or by straight-chain or branched alkyl having up to 6 carbon atoms, which, for its part, can be substituted by hydroxyl or by straight-chain or branched alkoxy having up to 4 carbon atoms,

D represents hydrogen, cycloalkyl having 4 to 12 carbon atoms or straight-chain or branched alkyl having up to 12 carbon atoms,

E represents the  $-\text{CO}-$  or  $-\text{CS}-$  group,

L represents an oxygen or sulphur atom or a group of the formula  $-\text{NR}^9$ ,

wherein

$R^9$  denotes hydrogen or straight-chain or branched alkyl having up to 6 carbon atoms, which is optionally substituted by hydroxyl or phenyl,

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$R^5$  represents phenyl or a 5- to 7-membered saturated or unsaturated heterocycle having up to 3 heteroatoms from the series consisting of S, N and/or O, the cycles optionally being substituted up to 3 times by identical or different nitro, carboxyl, halogen or cyano substituents or by straight-chain or branched alkenyl or alkoxy-carbonyl each having up to 6 carbon atoms or by straight-chain or branched alkyl having up to 6 carbon atoms, which is optionally substituted by hydroxyl, carboxyl or by straight-chain or branched alkoxy or alkoxycarbonyl each having up to 6 carbon atoms, and/or the cycles optionally being substituted by a group of the formula  $-\text{OR}^{10}$  or  $-\text{NR}^{11}\text{R}^{12}$ ,

wherein

$R^{10}$  denotes hydrogen or straight-chain or branched alkyl or alkenyl each having up to 6 carbon atoms,

$R^{11}$  and  $R^{12}$  are identical or different and denote phenyl, hydrogen or straight-chain or branched alkyl having up to 6 carbon atoms or straight-chain or branched acyl having up to 8 carbon atoms, which is optionally substituted by a group of the formula  $-\text{NR}^{13}\text{R}^{14}$ ,

wherein

$R^{13}$  and  $R^{14}$  are identical or different and denote hydrogen or straight-chain or branched acyl having up to 8 carbon atoms,

$R^6$  represents hydrogen, carboxyl or straight-chain or branched alkoxycarbonyl having up to 5 carbon atoms, or represents straight-chain or branched alkyl having up to 6 carbon atoms, which is optionally substituted by hydroxyl or by a group of the formula  $-\text{O}-\text{CO}-\text{R}^{15}$ ,

wherein

$R^{15}$  denotes phenyl which is optionally substituted up to 3 times by identical or different halogen or hydroxyl substituents or by straight-chain or branched alkyl having up to 5 carbon atoms, or straight-chain or branched alkyl or alkenyl each having up to 22 carbon atoms, each of which is optionally substituted by a group of the formula  $-\text{OR}^{16}$ ,

wherein

$R^{16}$  is hydrogen, benzyl, triphenylmethyl or straight-chain or branched acyl having up to 6 carbon atoms,

$R^7$  represents hydrogen or

$R^6$  and  $R^7$  together represent the group of the formula  $=\text{O}$ ,

if appropriate in an isomeric form, and their salts.

The cycloalkano-indole and -azaindole derivatives according to the invention can also be present in the form of their salts. In general, salts with organic or inorganic bases or acids may be mentioned here.

In the context of the present invention, physiologically acceptable salts are preferred. Physiologically acceptable salts of the compounds according to the invention can be salts of the substances according to the invention with mineral acids, carboxylic acids or sulphonic acids. Particularly preferred salts are, for example, those with hydrochloric acid, hydrobromic acid, sulphuric acid, phosphoric acid, methanesulphonic acid, ethanesulphonic acid, toluenesulphonic acid, benzenesulphonic acid, naphthalenedisulphonic acid, acetic acid, propionic acid, lactic acid, tartaric acid, citric acid, fumaric acid, maleic acid or benzoic acid.

Physiologically acceptable salts can also be metal or ammonium salts of the compounds according to the invention which have a free carboxyl group. Particularly preferred

salts are, for example, sodium, potassium, magnesium or calcium salts, and also ammonium salts which are derived from ammonia, or organic amines, such as, for example ethylamine, di- or triethylamine, di- or triethanolamine, dicyclohexylamine, dimethylaminoethanol, arginine, lysine, ethylenediamine or 2-phenylethylamine.

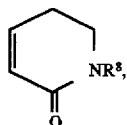
Including the double bond of parent structure, the cycloalkene radical ( $R^3/R^4$ ) in the context of the invention in general represents a 4- to 8-membered hydrocarbon radical, preferably a 5- to 8-membered hydrocarbon radical, for example a cyclobutene, cyclopentene, cyclohexene, cycloheptene or cyclooctene radical. The cyclopentene, cyclohexene, cyclooctene or cycloheptene radicals are preferred

Heterocycle ( $R^5$ ) in the context of the invention in general represents a saturated or unsaturated 5- to 7-membered heterocycle, preferably a 5- to 6-membered heterocycle, which can contain up to 3 heteroatoms from the series consisting of S, N and/or O. Examples which may be mentioned are: pyridyl, thienyl, furyl, pyrrolyl, thiazolyl, oxazolyl, imidazolyl, morpholinyl or piperidyl. Pyridyl and thienyl are preferred.

The compounds according to the invention can exist in stereoisomeric forms which either behave as image and mirror image (enantiomers), or do which do not behave as image and mirror image (diastereomers). The invention relates both to the enantiomers and diastereomers and their respective mixtures. These mixtures of the enantiomers and diastereomers can be separated in a known manner into the stereoisomerically uniform constituents.

Preferred compounds of the general formula (I) are those in which

$R^1$  and  $R^2$ , including the double bond connecting them, together form a phenyl or pyridyl ring or a ring of the formula



wherein

$R^3$  denotes hydrogen or straight-chain or branched alkyl having up to 3 carbon atoms,

$R^3$  and  $R^4$ , including the double bond connecting them, together form a phenyl ring or a cyclopentene, cyclohexene, cycloheptene, cyclooctene, oxocyclopentene, oxocyclohexene, oxocycloheptene or oxocyclooctene radical,

all ring systems mentioned under  $R^1/R^2$  and  $R^3/R^4$  optionally being substituted up to 2 times by identical or different fluorine, chlorine, bromine, trifluoromethyl, carboxyl or hydroxyl substituents, by straight-chain or branched alkoxy or alkoxy-carbonyl each having up to 4 carbon atoms or by straight-chain or branched alkyl having up to 4 carbon atoms, which, in turn, can be substituted by hydroxyl or by straight-chain or branched alkoxy having up to 3 carbon atoms,

D represents hydrogen, cyclobutyl, cyclopentyl, cyclohexyl, cycloheptyl, cyclooctyl or straight-chain or branched alkyl having up to 10 carbon atoms,

E represents the  $-\text{CO}-$  or  $-\text{CS}-$  group,

L represents an oxygen or sulphur atom or represents a group of the formula  $-\text{NR}^5$ ,

wherein

$R^9$  denotes hydrogen or straight-chain or branched alkyl having up to 5 carbon atoms, which is optionally substituted by hydroxyl or phenyl,

$R^5$  represents phenyl, pyridyl, furyl, thienyl or imidazolyl, each of which is optionally substituted up to 2 times by identical or different nitro, carboxyl, fluorine, chlorine, bromine or cyano substituents, by straight-chain or branched alkenyl or alkoxy carbonyl each having up to 4 carbon atoms or by straight-chain or branched alkyl having up to 5 carbon atoms, which is optionally substituted by hydroxyl, carboxyl or by straight-chain or branched alkoxy or alkoxy-carbonyl each having up to 5 carbon atoms,

and/or the cycles are optionally substituted by a group of the formula  $-\text{OR}^{10}$  or  $-\text{NR}^{11}\text{R}^{12}$ ,

wherein

$R^{10}$  denotes hydrogen or straight-chain or branched alkyl or alkenyl each having up to 4 carbon atoms,  $R^{11}$  and  $R^{12}$  are identical or different and denote phenyl, hydrogen or straight-chain or branched alkyl having up to 5 carbon atoms

or denote straight-chain or branched acyl having up to 6 carbon atoms, which is optionally substituted by a group of the formula  $-\text{NR}^{13}\text{R}^{14}$ ,

wherein

$R^{13}$  and  $R^{14}$  are identical or different and denote hydrogen or straight-chain or branched acyl having up to 6 carbon atoms,

$R^6$  represents hydrogen, carboxyl or straight-chain or branched alkoxy-carbonyl having up to 4 carbon atoms, or represents straight-chain or branched alkyl having up to 5 carbon atoms, which is optionally substituted by hydroxyl or by a group of the formula  $-\text{O}-\text{CO}-\text{R}^{15}$ ,

wherein

$R^{15}$  denotes phenyl which is optionally substituted up to 3 times by identical or different fluorine, chlorine, bromine or hydroxyl substituents or by straight-chain or branched alkyl having up to 4 carbon atoms, or straight-chain or branched alkyl or alkenyl each having up to 20 carbon atoms, each of which is optionally substituted by a group of the formula  $-\text{OR}^{16}$ ,

wherein

$R^{16}$  is hydrogen, benzyl, triphenylmethyl or straight-chain or branched acyl having up to 5 carbon atoms,

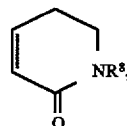
$R^7$  represents hydrogen or

$R^6$  and  $R^7$  together represent the group of the formula  $=\text{O}$ ,

if appropriate in an isomeric form, and their salts.

Particularly preferred compounds of the general formula (I) are those in which

$R^1$  and  $R^2$ , including the double bond connecting them, together form a phenyl or pyridyl ring or a ring of the formula



wherein

$R^8$  denotes hydrogen or methyl,

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R<sup>3</sup> and R<sup>4</sup>, including the double bond connecting them, together form a phenyl ring or a cyclopentene, cyclohexene, cycloheptene, cyclooctene, oxocyclopentene, oxocyclohexene, oxocycloheptene or oxocyclooctene radical,

all ring systems mentioned under R<sup>1</sup>/R<sup>2</sup> and R<sup>3</sup>/R<sup>4</sup> optionally being substituted up to 2 times by identical or different fluorine, chlorine, bromine, trifluoromethyl, carboxyl or hydroxyl substituents, by straight-chain or branched alkoxy or alkoxycarbonyl each having up to 3 carbon atoms or by straight-chain or branched alkyl having up to 3 carbon atoms, which, for its part, can be substituted by hydroxyl, methoxy or ethoxy,

D represents hydrogen, cyclopentyl, cyclohexyl, cycloheptyl, cyclooctyl or straight-chain or branched alkyl having up to 6 carbon atoms,

E represents the —CO— or —CS— group,

L represents an oxygen or sulphur atom or represents a group of the formula —NR<sup>9</sup>,

wherein

R<sup>9</sup> denotes hydrogen or straight-chain or branched alkyl having up to 4 carbon atoms, which is optionally substituted by hydroxyl or phenyl,

R<sup>5</sup> represents phenyl, pyridyl or thienyl, each of which is optionally substituted up to 2 times by identical or different nitro, carboxyl, fluorine, chlorine, bromine or cyano substituents, by straight-chain or branched alkyl or alkoxy carbonyl each having up to 3 carbon atoms or by straight-chain or branched alkyl having up to 4 carbon atoms, which is optionally substituted by hydroxyl, carboxyl or by straight-chain or branched alkoxy or alkoxycarbonyl each having up to 4 carbon atoms, and/or the cycles are optionally substituted by a group of the formula —OR<sup>10</sup> or —NR<sup>11</sup>R<sup>12</sup>,

wherein

R<sup>10</sup> denotes hydrogen or straight-chain or branched alkyl or alkenyl each having up to 3 carbon atoms,

R<sup>11</sup> and R<sup>12</sup> are identical or different and denote phenyl, hydrogen or straight-chain or branched alkyl having up to 4 carbon atoms

or denote straight-chain or branched acyl having up to 5 carbon atoms, which is optionally substituted by a group of the formula —NR<sup>13</sup>R<sup>14</sup>,

wherein

R<sup>13</sup> and R<sup>14</sup> are identical or different and denote hydrogen or straight-chain or branched acyl having up to 5 carbon atoms,

R<sup>6</sup> represents hydrogen, carboxyl or straight-chain or branched alkoxy carbonyl having up to 3 carbon atoms, or represents straight-chain or branched alkyl having up to 4 carbon atoms, which is optionally substituted by hydroxyl or by a group of the formula —CO—R<sup>15</sup>,

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wherein

R<sup>15</sup> denotes phenyl which is optionally substituted up to 3 times by identical or different straight-chain or branched alkyl having up to 3 carbon atoms, or denotes straight-chain or branched alkyl or alk-  
egyl each having up to 19 carbon atoms, each of which is optionally substituted by a group of the formula —OR<sup>16</sup>,

wherein

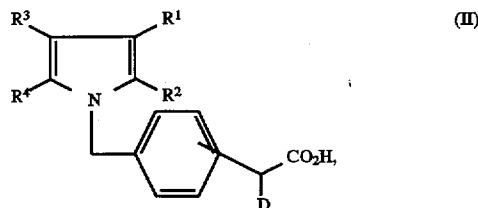
R<sup>16</sup> denotes hydrogen, benzyl, triphenylmethyl or straight-chain or branched acyl having up to 4 carbon atoms,

R<sup>7</sup> represents hydrogen or

R<sup>6</sup> and R<sup>7</sup> together represent the group of the formula =O,

if appropriate in an isomeric form, and their salts.

A process for the preparation of the compounds of the general formula (I) according to the invention has additionally been found, characterized in that carboxylic acids of the general formula (II)



in which

R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup> and D have the meaning indicated,

are amidated using compounds of the general formula (III)



in which

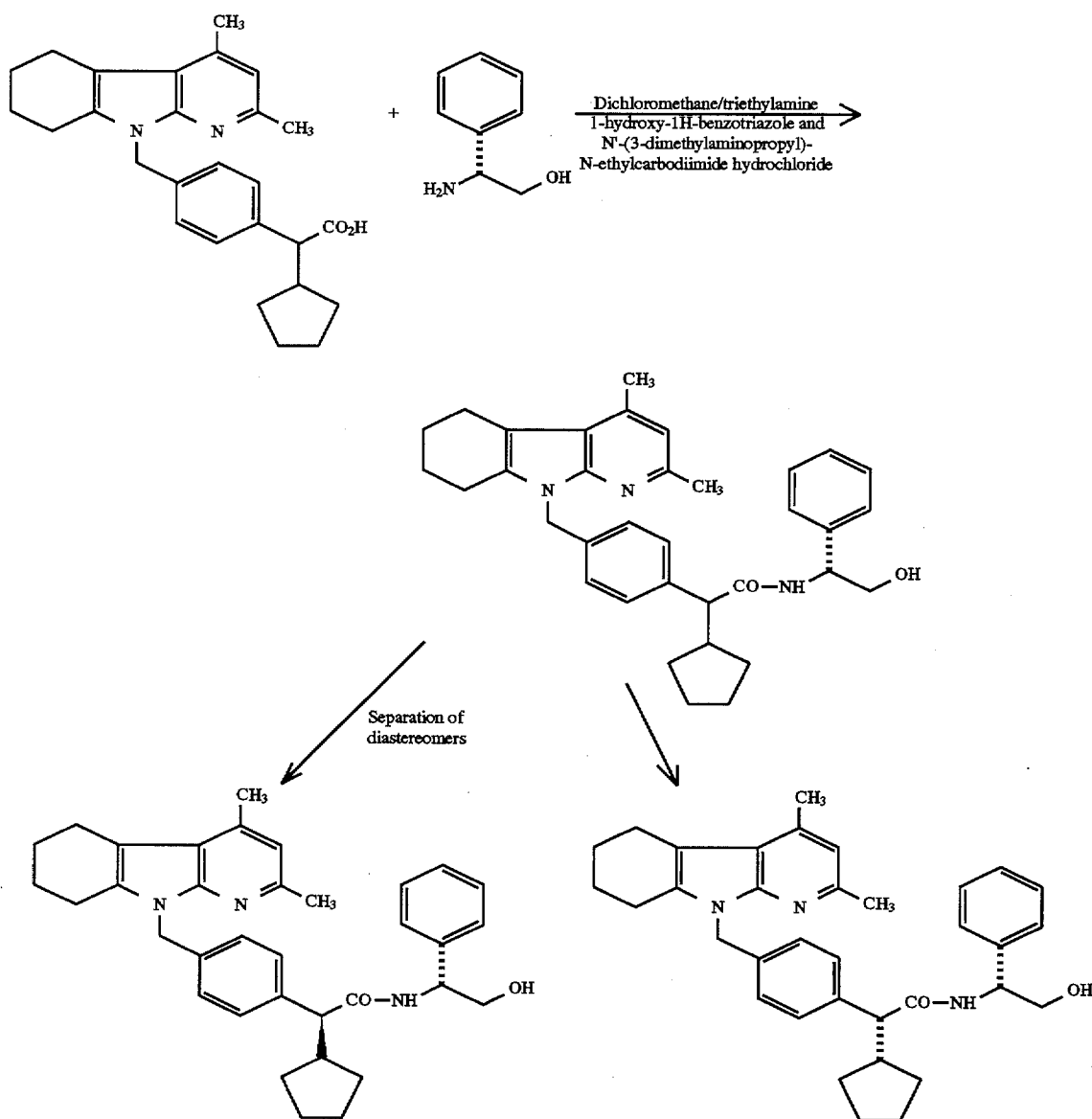
R<sup>5</sup> has the meaning indicated and

R<sup>17</sup> has the indicated meaning of R<sup>6</sup>, but does not represent carboxyl,

in an inert solvent and in the presence of bases and/or auxiliaries,

and, if appropriate, functional groups are varied by hydrolysis, esterification or reduction.

The process according to the invention can be illustrated by the following reaction scheme:



Suitable solvents for the amidation are in this case inert organic solvents which do not change under the reaction conditions. These include ethers, such as diethyl ether or tetrahydrofuran, halogenohydrocarbons such as dichloromethane, trichloromethane, tetrachloromethane, trichloroethane, tetrachloroethane, 1,2-dichloroethane or trichloroethylene, hydrocarbons such as benzene, xylene, toluene, hexane, cyclohexane or petroleum fractions, nitromethane, dimethylformamide, acetone, acetonitrile or hexamethylphosphoramide. It is also possible to employ mixtures of the solvents. Dichloromethane, tetrahydrofuran, acetone and dimethylformamide are particularly preferred.

Bases which can be employed for the process according to the invention are in general inorganic or organic bases. These preferably include alkali metal hydroxides such as, for example, sodium hydroxide or potassium hydroxide, alkaline earth metal hydroxides, such as, for example, barium hydroxide, alkali metal carbonates such as sodium carbonate or potassium carbonate, alkaline earth metal carbonates such as calcium carbonate, or alkali metal alkoxides such as

sodium or potassium methoxide, sodium or potassium ethoxide or potassium tert-butoxide, or organic amines (trialkyl(C<sub>1</sub>-C<sub>6</sub>)amines) such as triethylamine, or heterocycles such as 1,4-diazabicyclo[2.2.2]octane (DABCO), 1,8-diazabicyclo[5.4.0]undec-7-ene (DBU), pyridine, diaminopyridine, methylpiperidine or morpholine. It is also possible to employ alkali metals such as sodium and their hydrides such as sodium hydride as bases. Sodium and potassium carbonate and triethylamine are preferred.

The base is employed in an amount from 1 mol to 5 mol, preferably from 1 mol to 3 mol, relative to 1 mol of the compound of the general formula (II).

The reaction is in general carried out in a temperature range from 0° C. to 150° C., preferably from +20° C. to +110° C.

The reaction can be carried out at normal, increased or reduced pressure (e.g. 0.5 to 5 bar). In general, the reaction is carried out at normal pressure.

The amidation can optionally proceed via the activated stage of the acid halides, which can be prepared from the



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