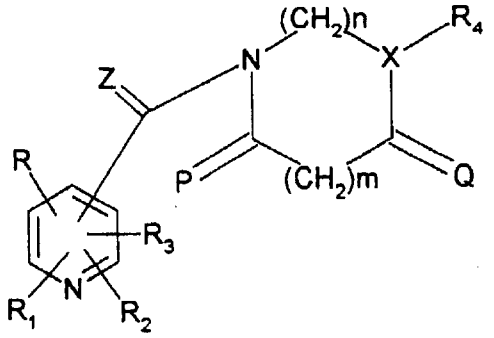
Novel heteroaryl derivatives and their use as medicaments

The invention relates to novel heteroaryl derivatives of the formula 1, to their preparation and to their use as medicaments, in particular for treating tumors.

According to one aspect of the invention, novel pyridine derivatives of the formula 1



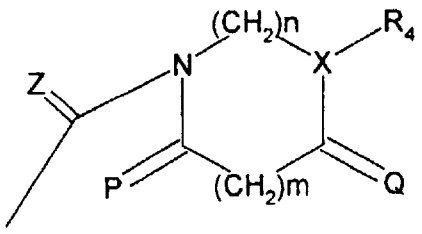
formula 1

in which

R, R_1, R_2, R_3 can be attached to any of the pyridine carbon atoms C_2 to C_6, are identical or different and independently of one another denote hydrogen, straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, straight-chain or branched (C_1-C_8)-alkylcarbonyl, preferably acetyl, straight-chain or branched (C_1-C_8)-alkoxy, halogen, aryl-(C_1-C_8)-alkoxy, preferably benzyloxy or phenylethyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, (C_1-C_8)-alkoxycarbonyl-amino, (C_1-C_6)-alkoxycarbonylamino-(C_1-C_8)-alkyl, cyano, straight-chain or branched cyano-(C_1-C_6)-alkyl, carboxyl, (C_1-C_8)-alkoxy-carbonyl, (C_1-C_4)-alkyl which is substituted by one or more fluorine atoms, preferably the trifluoromethyl group, carboxy-(C_1-C_8)-alkyl or (C_1-C_8)-alkoxycarbonyl-(C_1-C_6)-alkyl, $(C_2-$

C_6)-alkenyl, preferably allyl, (C_2-C_6)-alkynyl, preferably ethynyl or propargyl, straight-chain or branched cyano-(C_1-C_6)-alkyl, preferably cyanomethyl, aryl, where the aryl radical may be unsubstituted or mono- or polysubstituted by identical or different substituents from the group consisting of halogen, straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, carboxyl, straight-chain or branched (C_1-C_8)-alkoxycarbonyl, preferably tert-butoxycarbonyl, by trifluoromethyl, hydroxyl, straight-chain or branched (C_1-C_8)-alkoxy, preferably methoxy or ethoxy, benzyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, cyano, straight-chain or branched cyano-(C_1-C_6)-alkyl, where additionally R and R_1 and/or R_2 and R_3 may form a fused aromatic 6-membered ring with the pyridine ring forming a quinoline or acridine ring which for its part may be substituted at any C atom ring position by the radicals R, R_1, R_2 and R_3 having the meanings mentioned above;

Z is oxygen or sulfur, where the radical



substituted on the pyridine heterocycle may be attached to C atoms C_2-C_6 of the pyridine ring skeleton;

P, Q independently of one another represent oxygen or in each case two hydrogen atoms (i.e. -CH_2-);

X is nitrogen or C-R_5, where R_5 represents hydrogen or (C_1-C_6)-alkyl;

n,m independently of one another denotes an integer between 0-3, with the proviso that in the case n = 0, X denotes a CR_5R_6 group where R_5 and R_6 independently of one another represent hydrogen or (C_1-C_6)-alkyl and that the nitrogen atom adjacent to the C=Z group is substituted by a hydrogen atom or a (C-C_6)-alkyl group;

R_4 [lacuna] a straight-chain or branched (C_1-C_{20})-alkyl radical which may be saturated or unsaturated, with one to three double and/or triple bonds, and which may be unsubstituted or may optionally be substituted at the same or different C atoms by one, two or more aryl, heteroaryl, halogen, (C1-C6)-alkoxy, amino, mono-(C_1-C_4)-alkylamino or di-(C_1-C_4)-alkylamino; a (C_6-C_{14})-aryl radical, (C_6-C_{14})-aryl-(C_1-C_4)-alkyl radical or a (C_2-C_{10})-heteroaryl or (C_2-C_{10})-heteroaryl-(C_1-C_4)-alkyl radical which contains one or more heteroatoms selected from the group consisting of N, O and S, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identical or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and where the (C_6-C_{14})-aryl or (C_2-C_{10})-heteroaryl radical may be unsubstituted or mono- or polysubstituted by identical or different substituents from the group consisting of straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, halogen, cyano, (C_1-C_6)-alkoxycarbonylamino, (C_1-C_6)-alkoxy, carboxyl, (C_1-C_8)-alkoxycarbonyl, straight-chain or branched (C_1-C_6)-alkyl which is substituted by one or more fluorine atoms, preferably trifluoromethyl, hydroxyl, straight-chain or branched (C_1-C_8)-alkoxy, preferably methoxy or ethoxy, where

adjacent oxygen atoms may also be linked by (C_1-C_2)-alkylene groups, preferably by a methylene group, benzyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, aryl, which for its part may be unsubstituted or mono- or polysubstituted by identical or different substituents from the group consisting of straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, carboxyl, straight-chain or branched (C_1-C_8)-alkoxycarbonyl, by trifluoromethyl, hydroxyl, straight-chain or branched (C_1-C_8)-alkoxy, preferably methoxy or ethoxy, benzyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, cyano, straight-chain or branched cyano-(C_1-C_6)-alkyl;

and their structural isomers and stereoisomers, in particular tautomers, diastereomers and enantiomers, and their pharmaceutically acceptable salts, in particular acid addition salts, are provided.

Thus, for example, the compounds of the formula (1) according to the invention which have one or more centers of chirality and which are present as racemates can be separated by methods known per se into their optical isomers, i.e. enantiomers or diastereomers. The separation can be carried out by column separation on chiral phases or by recrystallization from an optically active solvent or using an optically active acid or base or by derivatization with an optically active reagent, such as, for example, an optically active alcohol, and subsequent removal of the radical.

Furthermore, the pyridine derivatives of the formula (1) according to the invention can be converted into their salts with inorganic or organic acids, in particular, for pharmaceutical use, into their physiologically acceptable salts. Acids which are

suitable for this purpose are, for example, hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, fumaric acid, succinic acid, lactic acid, citric acid, acetic acid, tartaric acid, malic acid, embonic acid, malonic acid, trifluoroacetic acid or maleic acid.

Moreover, the compounds of the formula (1) according to the invention can, if they contain a sufficiently acidic group, such as a carboxyl group, be converted, if desired, into their salts with inorganic or organic bases, in particular, for pharmaceutical use, into their physiologically acceptable salts. Bases which are suitable for this purpose are, for example, sodium hydroxide, potassium hydroxide, calcium hydroxide, lysine, cyclohexylamine, ethanolamine, diethanolamine and triethanolamine.

According to a preferred embodiment, pyridine derivatives of the formula 1 are provided in which R, R1, R2, R3, X, Z, P, Q, n and m have the meanings given above and

R_4 denotes a straight-chain or branched (C_1-C_{20})-alkyl radical which may be saturated or unsaturated, with one to three double and/or triple bonds, and which may be unsubstituted or optionally substituted on the same or different C atoms by one, two or more aryl, heteroaryl, halogen, (C1-C6)-alkoxy, amino, mono-(C_1-C_4)-alkylamino or di-(C_1-C_4)-alkylamino;

a phenyl ring or a naphthyl ring, each of which may be unsubstituted or mono- or polysubstituted by identical or different substituents from the group consisting of straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, halogen, cyano, (C_1-C_6)-alkoxycarbonylamino, (C_1-C_6)-alkoxy, carboxyl, (C_1-C_8)-alkoxycarbonyl, straight-chain or

branched (C_1-C_6)-alkyl which is substituted by one or more fluorine atoms, preferably trifluoromethyl, hydroxyl, straight-chain or branched (C_1-C_8)-alkoxy, preferably methoxy or ethoxy, where adjacent oxygen atoms may also be linked by (C_1-C_2)-alkylene groups, preferably a methylene group, benzyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, aryl, which for its part may be unsubstituted or mono- or polysubstituted by identical or different substituents from the group consisting of straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, carboxyl, straight-chain or branched (C_1-C_8)-alkoxycarbonyl, by trifluoromethyl, hydroxyl, straight-chain or branched (C_1-C_8)-alkoxy, preferably methoxy or ethoxy, benzyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, cyano, straight-chain or branched cyano-(C_1-C_6)-alkyl;

a 2-, 4-, 5- or 6-pyrimidinyl radical or 2-, 4-, 5- or 6-pyrimidinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 4-, 5- or 6-pyrimidinyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 3-, 4-, 5- or 6-pyridazinyl radical or 3-, 4-, 5- or 6-pyridazinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identical or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 3-, 4-, 5- or 6-pyridazinyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 5- or 6-pyrazinyl radical or 2-, 3-, 5- or 6-pyrazinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 3-, 5- or 6-pyrazinyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 3-, 4-, 5-, 6-, 7-, or 8-cinnolinyl radical or 3-, 4-, 5-, 6-, 7-, or 8-cinnolinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by

identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 3-, 4-, 5-, 6-, 7-, or 8-cinnolinyl radical may be unsubstituted or mono- to pentasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 4-, 5-, 6-, 7-, or 8-quinazolinyl radical or 2-, 4-, 5-, 6-, 7-, or 8-quinazolinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen and oxo (=O) and the or [sic] 2-, 4-, 5-, 6-, 7-, or 8-quinazolinyl radical may be unsubstituted or mono- to pentasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 5-, 6-, 7-, or 8-quinoxalinyl radical [lacuna] 2-, 3-, 5-, 6-, 7-, or 8-quinoxalinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the or [sic] 2-,

3-, 5-, 6-, 7-, or 8-quinoxalinyl radical may be unsubstituted or mono- to pentasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 4-, 5-, 6-, 7-, or 8-phthalazinyl radical or 1-, 4-, 5-, 6-, 7-, or 8-phthalazinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the or [sic] 1-, 4-, 5-, 6-, 7-, or 8-phthalazinyl radical may be unsubstituted or mono- to pentasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 4-, 5-, 6-, 7- or 8-quinolyl radical or 2-, 3-, 4-, 5-, 6-, 7 or 8-quinolyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 3-, 4-, 5-, 6-, 7- or 8-quinolyl radical may be unsubstituted or mono- to hexa-substituted by identical or different substituents

from the group consisting of hydrogen, (C_1-C_6)-alkyl, preferably methyl, particularly preferably 2-methyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 3-, 4-, 5-, 6-, 7- or 8-isoquinolyl radical or 1-, 3-, 4-, 5-, 6-, 7- or 8-isoquinolyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 3-, 4-, 5-, 6-, 7- or 8-isoquinolyl radical may be unsubstituted or mono- to hexa-substituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkyl-amino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxy-carbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 6-, 8- or 9-[9*H*]-purinyl radical or 2-, 6-, 8- or 9-[9*H*]-purinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 6-, 8- or 9-[9*H*]-purinyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino,

hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 6-, 7- or 8-[7*H*]-purinyl radical or 2-, 6-, 7- or 8-[7*H*]-purinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 6-, 7- or 8-[7*H*]-purinyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-acridinyl radical or 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-acridinyl-(C_1-C_4)-alkyl radical, where the (C_1-C6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-acridinyl radical may be unsubstituted or mono- to octasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoro-

methyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-phenanthridinyl radical or 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-phenanthridinyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-phenanthridinyl radical may be unsubstituted or mono- to octasubstituted by identical or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, (C_6-C_{10})-aryl-(C_1-C_6)-alkoxy, preferably benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 4-, 5- or 6-pyridyl radical where the 2-, 3-, 4-, 5- or 6-pyridyl radical may be unsubstituted or mono- to tetrasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 4-, 5- or 6-pyridinyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identical or different substituents from the group

consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 3-, 4-, 5- or 6-pyridinyl radical may be unsubstituted or mono- to tetrasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 4- or 5-thienyl radical or 2-, 3-, 4- or 5-thienyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 3-, 4- or 5-thienyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 4-, or 5-thiazolyl radical or 2-, 4-, or 5-thiazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 4-, or 5-thiazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of

hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 3-, 4-, or 5-isothiazolyl radical or 3-, 4-, or 5-isothiazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 3-, 4-, or 5-isothiazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 4-, 5-, 6-, or 7-benzothiazolyl radical or 2-, 4-, 5-, 6-, or 7-benzothiazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 4-, 5-, 6-, or 7-benzothiazolyl radical may be unsubstituted or mono- to tetrasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino

or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 2-, 4-, or 5-imidazolyl radical or 1-, 2-, 4-, or 5-imidazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 2-, 4-, or 5-imidazolyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 3-, 4-, or 5-pyrazolyl radical or 1-, 3-, 4- or 5-pyrazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 3-, 4- or 5-pyrazolyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 2-, 3-, 4-, or 5-pyrrolyl radical or 1-, 2-, 3-, 4-, or 5-pyrrolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 2-, 3-, 4- or 5-pyrrolyl radical may be unsubstituted or mono- to tetrasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 3-, or 5-[1.2.4]-triazolyl radical or 1-, 3-, or 5-[1.2.4]-triazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 3-, or 5-[1.2.4]-triazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 4-, or 5-[1.2.3]-triazolyl radical or 1-, 4-, or 5-[1.2.3]-triazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by

identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 4-, or 5-[1.2.3]-triazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1- or 5-[1*H*]-tetrazolyl radical or 1-, or 5-[1*H*]-tetrazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, or 5-[1*H*]-tetrazolyl radical may be unsubstituted or substituted by hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2- or 5-[2*H*]-tetrazoyl radical or 2- or 5-[2*H*]-tetrazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2- or 5-[2*H*]-tetrazolyl radical may be unsubstituted or substituted by hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-

alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 4-, or 6-[1.3.5]-triazinyl radical or 2-, 4-, or 6-[1.3.5]-triazinyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 4-, or 6-[1.3.5]-triazinyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 4-, or 5-oxazolyl radical or 2-, 4-, or 5-oxazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 4-, or 5-oxazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted

by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 3-, 4-, or 5-isoxazolyl radical or 3-, 4-, or 5-isoxazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 3-, 4-, or 5-isoxazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 2-, 3-, 4-, 5-, 6- or 7-indolyl radical or 1-, 2-, 3-, 4-, 5-, 6- or 7-indolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 2-, 3-, 4-, 5-, 6- or 7-indolyl radical may be unsubstituted or mono- to hexa-substituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl and the isomers, in particular tautomers, diastereomers and

enantiomers, and the pharmaceutically acceptable salts, in particular acid addition salts, thereof.

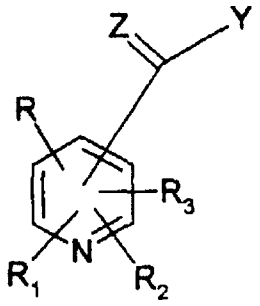
According to a further embodiment, pyridine derivatives of the formula (1) are provided which are characterized in that R, R_1, R_2, R_3, X, Z, P, Q, n and m have the meanings given above and R_4 represents phenyl which is unsubstituted or substituted by one to five identical or different (C_1-C_6)-alkoxy groups, where adjacent oxygen atoms may also be linked by (C_1-C_2)-alkylene groups.

According to a further embodiment, pyridine derivatives of the formula (1) are provided which are characterized in that R_4 has the meanings given above, R, R_1, R_2, R_3 each represent a hydrogen atom, Z represents an oxygen atom and X represents a nitrogen atom, P and Q each represent two hydrogen atoms (i.e. -CH2-), m is zero and n is the integer 2.

According to a further embodiment, pyridine derivatives of the formula (1) are provided which are characterized in that R, R_1, R_2, R_3, X, Z, P, Q, n and m have the meanings given above and R_4 represents 3,5-dimethoxyphenyl.

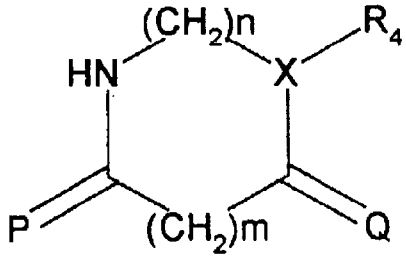
According to a further embodiment, pyridine derivatives of the formula (1) are provided which are characterized in that R, R_1, R_2, R_3 each represent a hydrogen atom, Z represents an oxygen atom, X represents a nitrogen atom, P and Q each represent two hydrogen atoms (i.e. -CH2-), m is zero, n represents the integer 2 and R_4 represents a 3,5-dimethoxyphenyl radical.

According to a further aspect of the invention, a process for preparing pyridine derivatives according to any of claims 1 to 6 is [lacuna], characterized in that a pyridine carboxylic acid of the formula (2)



formula 2

in which R, R1, R2, R3 have the meanings given above, Z denotes an oxygen or sulfur atom and Y represents a leaving group such as halogen, hydroxyl, (C1-C6)-alkoxy, preferably methoxy or ethoxy, -O-tosyl, -O-mesyl or imidazolyl, is reacted with an amine of the formula (3)



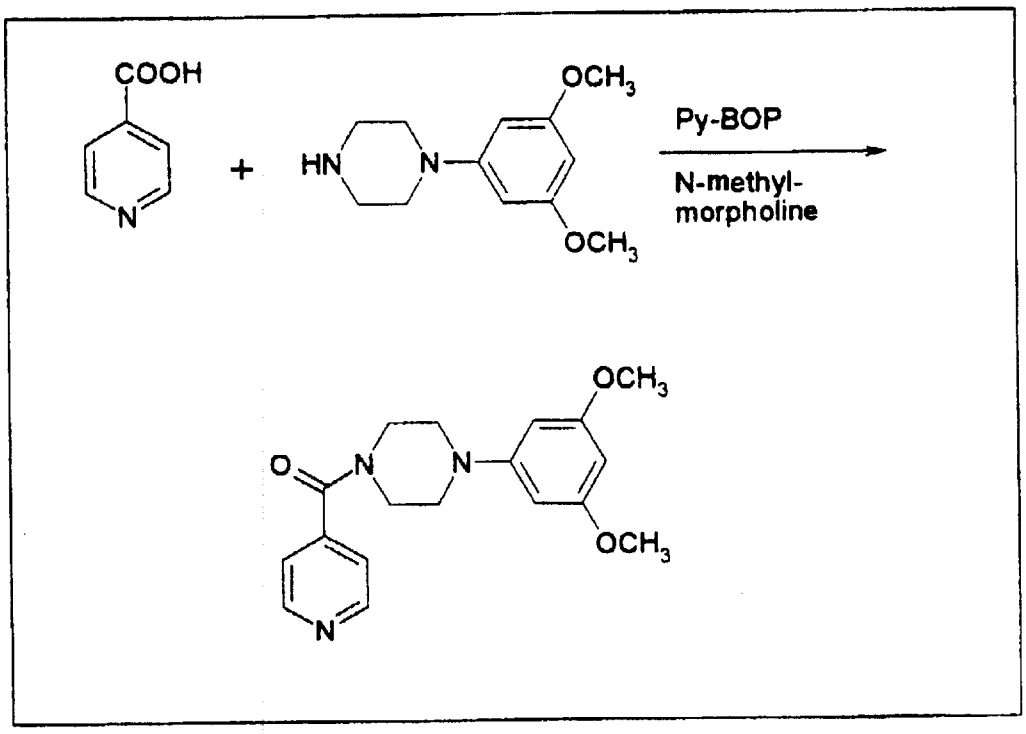
formula 3

in which R4, X, P, Q, m and n are as defined above, using, if appropriate, diluents and auxiliaries, and the desired pyridine derivative is formed.

Synthesis route:

The compounds of the formula 1 can be obtained according to Scheme 1 below:

Scheme 1



The starting materials (2) and (3) are either commercially available or can be prepared by procedures known per se. The starting materials (2) and (3) are useful intermediates for preparing the pyridine derivatives of the formula (1) according to the invention.

The solvents and auxiliaries to be used, if appropriate, and the reaction parameters to be used, such as reaction temperature and reaction time, are known from the literature or familiar to the person skilled in the art owing to his expert knowledge.

The pyridine derivatives of the formula (1) according to the invention are suitable as medicaments, in particular as antitumor agents, for treating mammals, in particular man, but also domestic animals such as horses, cattle, dogs, cats, hares, sheep, poultry and the like.

According to a further aspect of the invention, a method for controlling tumors in mammals, in particular man, is provided which is characterized in that at least one pyridine derivative of the formula (1) is administered to a mammal in an amount effective for the treatment of the tumor. The therapeutically effective dose of the pyridine derivative according to the invention in question which is to be administered for the treatment depends inter alia on the nature and the stage of the oncosis, the age and the sex of the patient, the type of administration and the duration of the treatment. Administration can take place orally, rectally, buccally (for example sublingually), parenterally (for example subcutaneously, intramuscularly, intradermally or intravenously), topically or transdermally.

According to a further aspect of the invention, medicaments for the treatment of tumors are provided which are characterized in that they comprise, as active ingredient, at least one pyridine derivative according to any of Claims 1 to 4 or a pharmaceutically acceptable salt thereof, if appropriate together with customary pharmaceutically acceptable auxiliaries, additives and carriers. These can be solid, semisolid, liquid or aerosol preparations. Suitable solid preparations are, for example, capsules, powders, granules, tablets. Suitable semisolid preparations are, for example, ointments, creams, gels, pastes, suspensions, oil-in-water and water-in-oil emulsions. Suitable liquid preparations are, for example, sterile aqueous preparations for parenteral administration which are isotonic with the blood of the patient.

The invention is to be illustrated in more detail by the example below, without being restricted to the example.

Working Example

1-(3,5-Dimethoxyphenyl)-4-(4-pyridyl-carbonyl)piperazine

5 g (45.8 mmol) of pyridine-4-carboxylic acid were suspended with stirring in 150 ml of DMF. With further stirring, 6.93 g (68.5 mmol) of N-methylmorpholine and then a solution of 35.6 g (68.5 mmol) of Py-BOP (1-benzotriazolyltripyrrolidinophosphonium hexafluorophosphate) and 10.18 g (45.8 mmol) of 1-(3,5-dimethoxyphenyl)piperazine in 60 ml of DMF were added to this mixture. The mixture was stirred at room temperature for 24 h, the DMF was distilled off under reduced pressure and the residue was purified on a silica gel column (Kieselgel 60, from Merck AG, Darmstadt) using the mobile phase dichloromethane/methanol/25 percent ammonia (90:10:1 v/v/v).

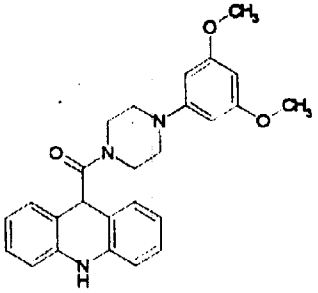
Yield: 12.34 g (82.3% of theory)
m.p.: 94-96°C

1. Antiproliferative action in various tumor cell lines

In a proliferation test, the antiproliferative activity of the substance D-43411 was examined using established tumor cell lines. In the test used, the cellular dehydrogenation activity is determined, which makes it possible to determine the vitality of the cell and, indirectly, the cell count. The cell lines used are the human cervical carcinoma cell lines KB/HeLa (ATCC/CCL17), the murine lymphocyte leukaemia L1210 (ATCC CCL-219), the human breast adenocarcinoma line MCF7 /ATCC HTB22) and the ovary adenocarcinoma line SKOV-3 (ATCC HTB77). These are established cell lines which are very well characterized and were obtained from ATCC and cultured.

The results shown in Tab. 1 demonstrate the highly potent antiproliferative action of D-43411 in the cell lines SKOV-3, L-1210 and HeLa/KB. Owing to the particularly slow growth of the MCF7 line, the effect of D-43411 in the test period of 48 h is only small (18% inhibition at 3.16 µg/ml; thus stated as >3.16).

Tab. 1 In-vitro cytotoxicity in tumor cell lines (values determined from 5 substance concentrations)

		XTT - Assay IC^{50} [µg/ml]				
D number	Structure	MW	SKOV-3	L1210	KB/HeLa	MCF7
D-43411		429	<0.0003	<0.0003	<0.0003	>3.16

2. Method

XTT Test for cellular dehydrogenase activity

The adherently growing tumor cell lines HeLa/KB, SKOV-3 and MCF7 and the L1210 leukaemia line, which grows in suspension, were cultivated under standard conditions in an incubator with gas inlet at 37°C, 5% CO_2 and 95% atmospheric humidity. On Test Day 1, the adherent cells are detached using trypsine/EDTA and pelleted by centrifugation. The cell pellet is then resuspended in RPMI culture medium at the appropriate cell count and transferred to a 96-well microtitre plate. The plates are then cultivated overnight in the incubator with gas inlet. The test substances are made up as stock solutions in DMSO and, on Test Day 2, diluted with culture medium to the desired concentrations. The

substances in the culture medium are then added to the cells and incubated in the incubator with gas inlet for 45 h. Cells which have not been treated with test substance serve as control.

For the XTT assay, 1 mg/ml of XTT (sodium 3'-[1-(phenylaminocarbonyl)-3,4-tetrazolium]-bis(4-methoxy-6-nitro)benzenesulfonic acid) is dissolved in RPMI-1640 medium without Phenol Red. Additionally, a 0.383 mg/ml solution of PMS (N-methyldibenzopyrazine methyl sulfate) in phosphate-buffered saline (PBS) is prepared. On Test Day 4, 75 µl/well of the XTT-PMS mixture are pipetted onto the cell plates, which by now have been incubated with the test substances for 45 h. To this end, the XTT solution is mixed with the PMS solution in a ratio of 50:1 (v/v) shortly before use. The cell plates are then incubated in the incubator with gas inlet for a further 3 h, and the optical density (OD_{490nm}) is determined in a photometer.

Using the OD_{490nm} obtained, the inhibition in percent relative to the control is calculated. The antiproliferative activity is estimated using regression analysis.

Example I

Tablet containing 50 mg of active compound

Composition:

(1)	Active compound	50.0 mg
(2)	Lactose	98.0 mg
(3)	Maize starch	50.0 mg
(4)	Polyvinylpyrrolidone	15.0 mg
(5)	Magnesium stearate	2.0 mg
Total:		215.0 mg

Preparation:

(1), (2) and (3) are mixed and granulated with an aqueous solution of (4). The dried granules are admixed with (5). This mixture is tabletted.

Example II

Capsule containing 50 mg of active compound

Composition:

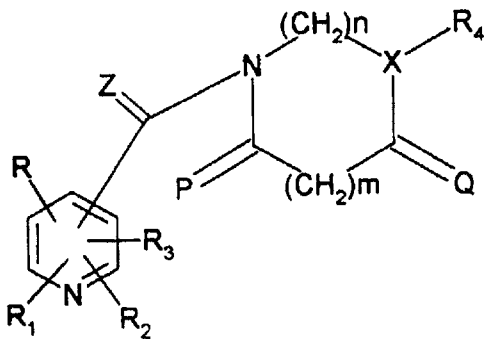
(1)	Active compound	50.0 mg
(2)	Maize starch, dried	58.0 mg
(3)	Lactose powder	50.0 mg
(4)	Magnesium stearate	2.0 mg
Total:		160.0 mg

Preparation:

(1) is ground with (3). This ground material is added with vigorous mixing to the mixture of (2) and (4). This powder mixture is, on a capsule filling machine, filled into hard gelatine capsules size 3.

Patent Claims

1. Pyridine derivatives according to the formula 1



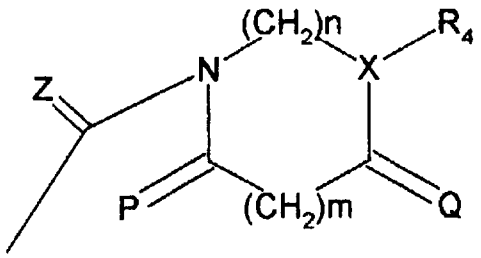
formula 1

in which

R, R_1, R_2, R_3 can be attached to any of the pyridine carbon atoms C_2 to C_8, are identical or different and independently of one another denote hydrogen, straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, straight-chain or branched (C_1-C_8)-alkylcarbonyl, preferably acetyl, straight-chain or branched (C_1-C_8)-alkoxy, halogen, aryl-(C_1-C_8)-alkoxy, preferably benzyloxy or phenylethyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, (C_1-C_8)-alkoxycarbonyl-amino, (C_1-C_6)-alkoxycarbonylamino-(C_1-C_8)-alkyl, cyano, straight-chain or branched cyano-(C_1-C_6)-alkyl, carboxyl, (C_1-C_8)-alkoxy-carbonyl, (C_1-C_4)-alkyl which is substituted by one or more fluorine atoms, preferably the trifluoromethyl group, carboxy-(C_1-C_8)-alkyl or (C_1-C_8)-alkoxycarbonyl-(C_1-C_6)-alkyl, (C_2-C_6)-alkenyl, preferably allyl, (C_2-C_6)-alkynyl, preferably ethynyl or propargyl, straight-chain or branched cyano-(C_1-C_6)-alkyl, preferably cyanomethyl, aryl, where the aryl radical may be unsubstituted or mono- or polysubstituted by identical or

different substituents from the group consisting of halogen, straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, carboxyl, straight-chain or branched (C_1-C_8)-alkoxycarbonyl, preferably tert-butoxycarbonyl, by trifluoromethyl, hydroxyl, straight-chain or branched (C_1-C_8)-alkoxy, preferably methoxy or ethoxy, benzyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, cyano, straight-chain or branched cyano-(C_1-C_6)-alkyl, where additionally R and R_1 and/or R_2 and R_3 may form a fused aromatic 6-membered ring with the pyridine ring forming a quinoline or acridine ring which for its part may be substituted at any C atom ring position by the radicals R, R_1, R_2 and R_3 having the meanings mentioned above;

Z is oxygen or sulfur, where the radical



substituted on the pyridine heterocycle may be attached to C atoms C_2-C_6 of the pyridine ring skeleton;

P,Q indepedently of one another represent oxygen or in each case two hydrogen atoms (i.e. $-CH_2-$);

X is nitrogen or C-R_5, where R_5 represents hydrogen or (C_1-C_6)-alkyl;

n,m independently of one another denotes an integer between 0-3, with the proviso that in the case n = 0, X denotes a CR_5R_6 group where R_5 and R_6 independently of one another represent hydrogen or (C_1-C_6)-alkyl and that the nitrogen atom adjacent to the C=Z group is substituted by a hydrogen atom or a $(C-C_6)$-alkyl group;

R_4 [lacuna] a straight-chain or branched (C_1-C_{20})-alkyl radical which may be saturated or unsaturated, with one to three double and/or triple bonds, and which may be unsubstituted or may optionally be substituted at the same or different C atoms by one, two or more aryl, heteroaryl, halogen, (C1-C6)-alkoxy, amino, mono-(C_1-C_4)-alkylamino or di-(C_1-C_4)-alkylamino; a (C_6-C_{14})-aryl radical, (C_6-C_{14})-aryl-(C_1-C_4)-alkyl radical or a (C_2-C_{10})-heteroaryl or (C_2-C_{10})-heteroaryl-(C_1-C_4)-alkyl radical which contains one or more heteroatoms selected from the group consisting of N, O and S, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identical or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and where the (C_6-C_{14})-aryl or (C_2-C_{10})-heteroaryl radical may be unsubstituted or mono- or polysubstituted by identical or different substituents from the group consisting of straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, halogen, cyano, (C_1-C_6)-alkoxycarbonylamino, (C_1-C_6)-alkoxy, carboxyl, (C_1-C_8)-alkoxycarbonyl, straight-chain or branched (C_1-C_6)-alkyl which is substituted by one or more fluorine atoms, preferably trifluoromethyl, hydroxyl, straight-chain or branched (C_1-C_8)-alkoxy, preferably methoxy or ethoxy, where adjacent oxygen atoms may also be linked by (C_1-C_2)-alkylene groups, preferably by a methylene group, benzyloxy, nitro, amino, mono-(C_1-C_4)-

alkylamino, di-(C_1-C_4)-alkylamino, aryl, which for its part may be unsubstituted or mono- or polysubstituted by identical or different substituents from the group consisting of straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, carboxyl, straight-chain or branched (C_1-C_8)-alkoxycarbonyl, by trifluoromethyl, hydroxyl, straight-chain or branched (C_1-C_8)-alkoxy, preferably methoxy or ethoxy, benzyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, cyano, straight-chain or branched cyano-(C_1-C_6)-alkyl;

and their structural isomers and stereoisomers, in particular tautomers, diastereomers and enantiomers, and their pharmaceutically acceptable salts, in particular acid addition salts.

2. Pyridine derivatives of the formula 1 according to Claim 1, characterized in that R, R1, R2, R3, P, Q, X, Z, n and m have the meanings given in Claim 1 and

R4 denotes a straight-chain or branched (C_1-C_{20})-alkyl radical which may be saturated or unsaturated, with one to three double and/or triple bonds, and which may be unsubstituted or optionally substituted on the same or different C atoms by one, two or more aryl, heteroaryl, halogen, (C1-C6)-alkoxy, amino, mono-(C_1-C_4)-alkylamino or di-(C_1-C_4)-alkylamino;

a phenyl ring or a naphthyl ring, each of which may be unsubstituted or mono- or poly-substituted by identical or different substituents from the group consisting of straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, halogen, cyano, (C_1-C_6)-alkoxycarbonylamino, (C_1-C_6)-alkoxy, carboxyl,

(C_1-C_8)-alkoxycarbonyl, straight-chain or branched (C_1-C_6)-alkyl which is substituted by one or more fluorine atoms, preferably trifluoromethyl, hydroxyl, straight-chain or branched (C_1-C_8)-alkoxy, preferably methoxy or ethoxy, benzyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, aryl, which for its part may be unsubstituted or mono- or polysubstituted by identical or different substituents from the group consisting of straight-chain or branched (C_1-C_8)-alkyl, (C_3-C_7)-cycloalkyl, carboxyl, straight-chain or branched (C_1-C_8)-alkoxycarbonyl, by trifluoromethyl, hydroxyl, straight-chain or branched (C_1-C_8)-alkoxy, preferably methoxy or ethoxy, benzyloxy, nitro, amino, mono-(C_1-C_4)-alkylamino, di-(C_1-C_4)-alkylamino, cyano, straight-chain or branched cyano-(C_1-C_6)-alkyl;

a 2-, 4-, 5- or 6-pyrimidinyl radical or 2-, 4-, 5- or 6-pyrimidinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 4-, 5- or 6-pyrimidinyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 3-, 4-, 5- or 6-pyridazinyl radical or 3-, 4-, 5- or 6-pyridazinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identical or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 3-, 4-, 5- or 6-pyridazinyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonyl-amino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 5- or 6-pyrazinyl radical or 2-, 3-, 5- or 6-pyrazinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 3-, 5- or 6-pyrazinyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 3-, 4-, 5-, 6-, 7-, or 8-cinnolinyl radical or 3-, 4-, 5-, 6-, 7-, or 8-cinnolinyl-(C_1-C_4)-

alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 3-, 4-, 5-, 6-, 7-, or 8-cinnolinyl radical may be unsubstituted or mono- to pentasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 4-, 5-, 6-, 7-, or 8-quinazolinyl radical or 2-, 4-, 5-, 6-, 7-, or 8-quinazolinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen and oxo (=O) and the or [sic] 2-, 4-, 5-, 6-, 7-, or 8-quinazolinyl radical may be unsubstituted or mono- to pentasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 5-, 6-, 7-, or 8-quinoxalinyl radical [lacuna] 2-, 3-, 5-, 6-, 7-, or 8-quinoxalinyl-

(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the or [sic] 2-, 3-, 5-, 6-, 7-, or 8-quinoxalinyl radical may be unsubstituted or mono- to pentasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonyl-amino or (C_1-C_6)-alkyl which is mono- or poly-substituted by fluorine, preferably trifluoro-methyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 4-, 5-, 6-, 7-, or 8-phthalazinyl radical or 1-, 4-, 5-, 6-, 7-, or 8-phthalazinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- [lacuna] polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the or [sic] 1-, 4-, 5-, 6-, 7-, or 8-phthalazinyl radical may be unsubstituted or mono- to pentasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonyl-amino or (C_1-C_6)-alkyl which is mono- or poly-substituted by fluorine, preferably trifluoro-methyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 4-, 5-, 6-, 7- or 8-quinolyl radical or 2-, 3-, 4-, 5-, 6-, 7 or 8-quinolyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 3-, 4-, 5-, 6-, 7- or 8-quinolyl radical may be unsubstituted or mono- to hexasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, preferably methyl, particularly preferably 2-methyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 3-, 4-, 5-, 6-, 7- or 8-isoquinolyl radical or 1-, 3-, 4-, 5-, 6-, 7- or 8-isoquinolyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 3-, 4-, 5-, 6-, 7- or 8-isoquinolyl radical may be unsubstituted or mono- to hexasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 6-, 8- or 9-[9*H*]-purinyl radical or 2-, 6-, 8- or 9-[9*H*]-purinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 6-, 8- or 9-[9*H*]-purinyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonyl-amino or (C_1-C_6)-alkyl which is mono- or poly-substituted by fluorine, preferably trifluoro-methyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 6-, 7- or 8-[7*H*]-purinyl radical or 2-, 6-, 7- or 8-[7*H*]-purinyl-(C_1-C_4)-alkyl radical, where the (C_1-C_4)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 6-, 7- or 8-[7*H*]-purinyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonyl-amino or (C_1-C_6)-alkyl which is mono- or poly-substituted by fluorine, preferably trifluoro-methyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-acridinyl radical or 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-acridinyl-(C_1-C_4)-alkyl radical, where the (C1-C6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-acridinyl radical may be unsubstituted or mono- to octasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-phenanthridinyl radical or 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-phenanthridinyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 2-, 3-, 4-, 5-, 6-, 7-, 8- or 9-phenanthridinyl radical may be unsubstituted or mono- to octasubstituted by identical or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, (C_6-C_{10})-aryl-(C_1-C_6)-alkoxy, preferably benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine,

preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 4-, 5- or 6-pyridyl radical where the 2-, 3-, 4-, 5- or 6-pyridinyl [sic] radical may be unsubstituted or mono- to tetrasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 4-, 5- or 6-pyridyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identical or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 3-, 4-, 5- or 6-pyridinyl radical may be unsubstituted or mono- to tetrasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 3-, 4- or 5-thienyl radical or 2-, 3-, 4- or 5-thienyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group

consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 3-, 4- or 5-thienyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 4-, or 5-thiazolyl radical or 2-, 4-, or 5-thiazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 4-, or 5-thiazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 3-, 4-, or 5-isothiazolyl radical or 3-, 4-, or 5-isothiazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 3-, 4-, or 5-isothiazolyl radical may be unsubstituted or mono- or disubstituted

by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 4-, 5-, 6-, or 7-benzothiazolyl radical or 2-, 4-, 5-, 6-, or 7-benzothiazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 4-, 5-, 6-, or 7-benzothiazolyl radical may be unsubstituted or mono- to tetrasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 2-, 4-, or 5-imidazolyl radical or 1-, 2-, 4-, or 5-imidazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 2-, 4-, or 5-imidazolyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of

hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 3-, 4-, or 5-pyrazolyl radical or 1-, 3-, 4- or 5-pyrazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 3-, 4- or 5-pyrazolyl radical may be unsubstituted or mono- to trisubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 2-, 3-, 4-, or 5-pyrrolyl radical or 1-, 2-, 3-, 4-, or 5-pyrrolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 2-, 3-, 4- or 5-pyrrolyl radical may be unsubstituted or mono- to tetrasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino,

hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 3-, or 5-[1.2.4]-triazolyl radical or 1-, 3-, or 5-[1.2.4]-triazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 3-, or 5-[1.2.4]-triazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 4-, or 5-[1.2.3]-triazolyl radical or 1-, 4-, or 5-[1.2.3]-triazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 4-, or 5-[1.2.3]-triazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl,

(C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1- or 5-[1*H*]-tetrazolyl radical or 1-, or 5-[1*H*]-tetrazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, or 5-[1*H*]-tetrazolyl radical may be unsubstituted or substituted by hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2- or 5-[2*H*]-tetrazoyl radical or 2- or 5-[2*H*]-tetrazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2- or 5-[2*H*]-tetrazolyl radical may be unsubstituted or substituted by hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 4-, or 6-[1.3.5]-triazinyl radical or 2-, 4-, or 6-[1.3.5]-triazinyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 4-, or 6-[1.3.5]-triazinyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 2-, 4-, or 5-oxazolyl radical or 2-, 4-, or 5-oxazolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 2-, 4-, or 5-oxazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 3-, 4-, or 5-isoxazolyl radical or 3-, 4-, or 5-isoxazolyl-(C_1-C_6)-alkyl radical, where the

(C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 3-, 4-, or 5-isoxazolyl radical may be unsubstituted or mono- or disubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl;

a 1-, 2-, 3-, 4-, 5-, 6- or 7-indolyl radical or 1-, 2-, 3-, 4-, 5-, 6- or 7-indolyl-(C_1-C_6)-alkyl radical, where the (C_1-C_6)-alkyl radical may be unsubstituted or mono- or polysubstituted by identidal or different substituents from the group consisting of (C_1-C_6)-alkyl, halogen and oxo (=O) and the 1-, 2-, 3-, 4-, 5-, 6- or 7-indolyl radical may be unsubstituted or mono- to hexasubstituted by identical or different substituents from the group consisting of hydrogen, (C_1-C_6)-alkyl, halogen, nitro, amino, mono-(C_1-C_6)-alkylamino, di-(C_1-C_6)-alkylamino, hydroxyl, (C_1-C_6)-alkoxy, benzyloxy, carboxyl, (C_1-C_6)-alkoxycarbonyl, (C_1-C_6)-alkoxycarbonylamino or (C_1-C_6)-alkyl which is mono- or polysubstituted by fluorine, preferably trifluoromethyl, (C_6-C_{10})-aryl and (C_6-C_{10})-aryl-(C_1-C_6)-alkyl.

3. Pyridine derivatives according to Claim 1 or 2, characterized in that R, R_1, R_2, R_3, P, Q, X, Z, n and m have the meanings given above and R4 represents phenyl which is unsubstituted or

substituted by one to five identical or different (C_1-C_6)-alkoxy groups, where adjacent oxygen atoms may also be linked by (C_1-C_2)-alkylene groups.

4. Pyridine derivatives according to any of Claims 1 to 3, characterized in that R, R_1, R_2, R_3, P, Q, X, Z, n and m have the meanings given above and R_4 represents 3,5-dimethoxyphenyl.

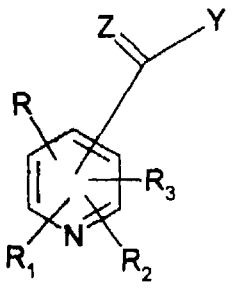
5. Pyridine derivatives according to any of the preceding claims, characterized in that R_4 has the meanings given above, R, R_1, R_2, R_3 each represent a hydrogen atom, Z represents an oxygen atom and X represents a nitrogen atom, P and Q each represent two hydrogen atoms (i.e. -CH2-), m is zero, n represents the integer 2.

6. Pyridine derivative according to any of Claims 1 to 5, characterized in that R, R_1, R_2, R_3 each represent a hydrogen atom, Z represents an oxygen atom, X represents a nitrogen atom, P and Q each represent two hydrogen atoms (i.e. -CH2-), m is zero, n represents the integer 2, and R_4 represents a 3,5-dimethoxyphenyl radical.

7. Pyridine derivatives according to any of Claims 1 to 6 for use as medicaments.

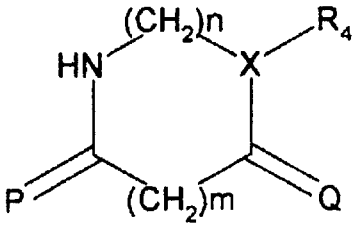
8. Use of the pyridine derivatives according to any of Claims 1 to 6 for preparing a medicament for treating tumors in mammals.

9. Process for preparing pyridine derivatives according to any of Claims 1 to 6, characterized in that a pyridine carboxylic acid of the formula (2)



formula 2

in which R, R1, R2, R3 have the meanings given above, Z denotes an oxygen or sulfur atom and Y represents a leaving group such as halogen, hydroxyl, (C1-C6)-alkoxy, preferably methoxy or ethoxy, -O-tosyl, -O-mesyl or imidazolyl, is reacted with an amine of the formula (3)



formula 3

in which R4, X, P, Q, m and n have the meanings given above, if appropriate using diluents and auxiliaries, and the desired pyridine derivative is formed.

10. Method for treating tumors in mammals, characterized in that at least one pyridine derivative according to any of Claims 1 to 6 is administered to the mammal in a dose effective for tumor treatment.

11. Medicament, characterized in that it comprises, as active ingredient, at least one pyridine derivative according to any of Claims 1 to 6, if appropriate together with customary

pharmaceutically acceptable auxiliaries, additives and carriers.