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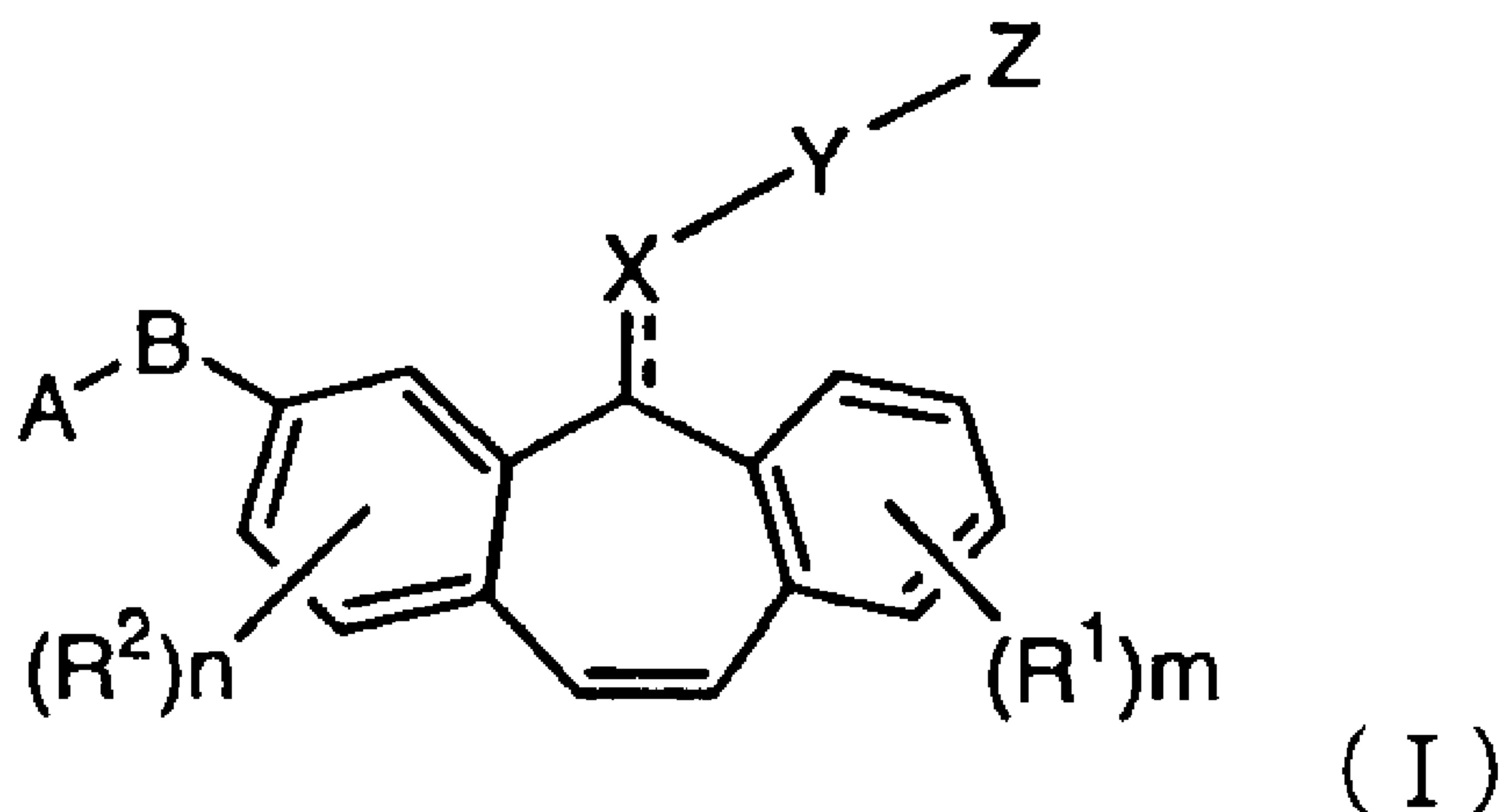
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(54) Titre : COMPOSE DE DIBENZOCYCLOHEPTENE

(54) Title: DIBENZOCYCLOHEPTENE COMPOUND



(57) **Abrégé/Abstract:**

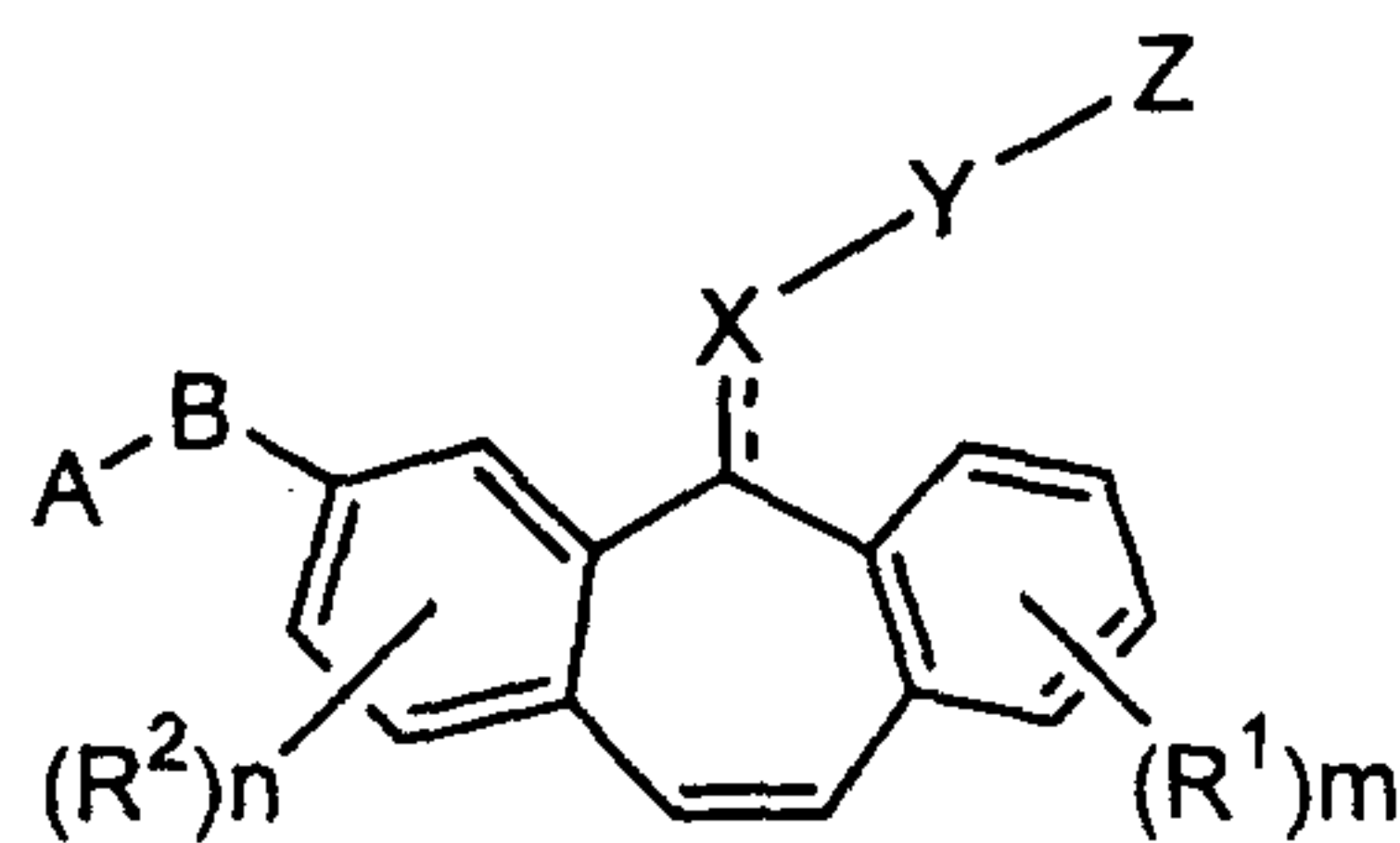
A dibenzocycloheptene compound represented by the general formula (I): (I) (wherein R¹ represents hydrogen, halogeno, etc.; R² represents hydrogen, halogeno, etc.; A represents, e.g., a 5- or 6-membered heterocyclic aromatic group containing one to three heteroatoms selected from the group consisting of nitrogen, oxygen, and sulfur, the heterocyclic aromatic group optionally having halogeno, nitro, etc. as a substituent; B represents -CH=CH-, -CH₂O-, etc.; Y represents C₁₋₁₀ alkylene optionally having halogeno, etc. as a substituent; Z represents optionally protected carboxyl, etc.; m is an integer of 1 to 4; n is an integer of 1 to 3; and indicates a single bond or double bond) or a pharmacologically acceptable salt of the compound; and a medicinal composition which contains the compound or salt as the active ingredient. The composition has not only strong leucotriene D₄ antagonism but leucotriene C₄ antagonism and leucotriene E₄ antagonism. It is useful as an antasthmatic, antiallergic, or antiphlogistic.



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Abstract

The present invention discloses a dibenzocycloheptene compound represented by the formula (I):



(I)

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10

15

wherein R^1 : hydrogen atom, halogen atom, etc., R^2 : hydrogen atom, halogen atom, etc., A: 5-membered or 6-membered heteroaromatic ring group containing 1 to 3 hetero atom(s) selected from the group consisting of nitrogen atom, oxygen atom and sulfur atom, and the heteroaromatic ring group, etc. may have halogen atom, nitrogen atom, etc. as substituent(s), B: formula; $-\text{CH}=\text{CH}-$, formula; $-\text{CH}_2\text{O}-$, etc., Y: C_1 - C_{10} alkylene group which may have halogen atom, etc. as substituent(s), etc., Z: carboxyl group which may be protected, etc., m: an integer of 1 to 4, n: an integer of 1 to 3,

20

----- represents a single bond or a double bond, or a pharmaceutically acceptable salt thereof and a medical composition containing the same as an effective ingredient which has leukotriene C_4 antagonistic action and leukotriene E_4 antagonistic action in addition to potent leukotriene D_4 antagonistic action, and useful as antiasthmatic agent, antiallergic agent and anti-inflammatory agent.

Specification

DIBENZOCYCLOHEPTENE COMPOUND

5 Technical field

The present invention relates to a dibenzocycloheptene compound which has, in addition to potent leukotriene D₄ antagonistic action, leukotriene C₄ antagonistic action and leukotriene E₄ antagonistic action, and useful
10 as an antiazma agent, an antiallergic agent and anti-inflammatory agent, or a pharmaceutically acceptable salt thereof.

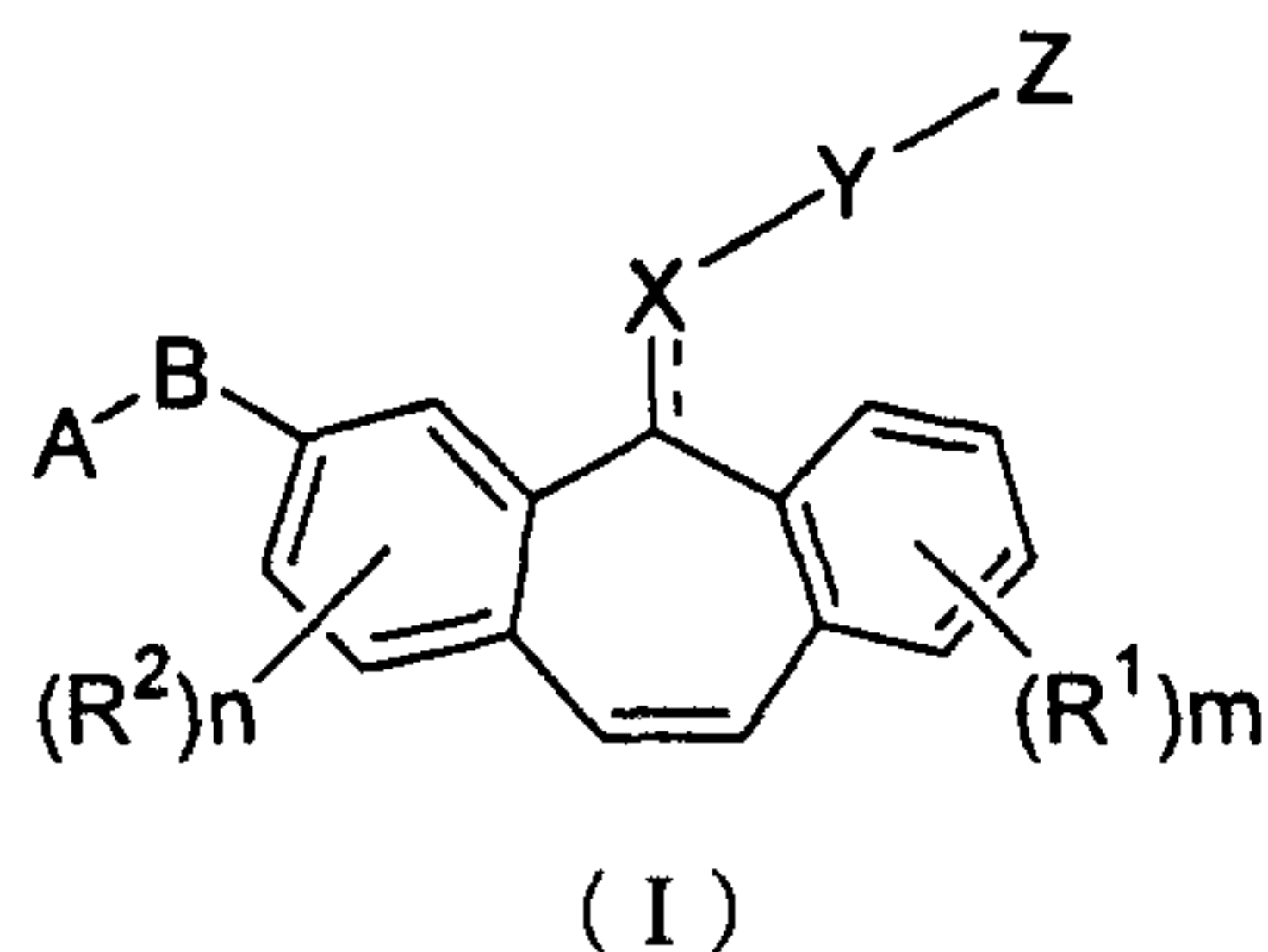
Background art

15 As a compound having leukotriene D₄ antagonistic action similarly in the present invention and having a structure similar to the compound of the present invention, there has been known, for example, a compound disclosed in WO94/19345, and as a compound having a structure partially
20 similar to the same, there have been known 5-[3-[3-(2-quinolinylmethoxy)phenoxy]propyl]-1H-tetrazole (RG7152; J. Med. Chem., 33, 1186 (1990)), 5-[[2-[[4-(2-quinolinylmethoxy)phenoxy]methyl]phenyl]methyl]-1H-tetrazole (RG12525; J. Med. Chem., 33, 1194 (1990)), and a compound
25 disclosed in WO95/18107.

In the present invention, as a result of research for long years about syntheses of compounds having potent leukotriene D₄ antagonistic action, as well as having antagonistic actions to leukotriene C₄ and leukotriene E₄
30 and their pharmaceutical effects, the inventors have found that novel dibenzocycloheptene compounds have excellent leukotriene D₄ antagonistic action, as well as having leukotriene C₄ and leukotriene E₄ antagonistic action with good balance, and have high safety, excellent oral absorb-
35 ability and durability of the action to accomplish the present invention.

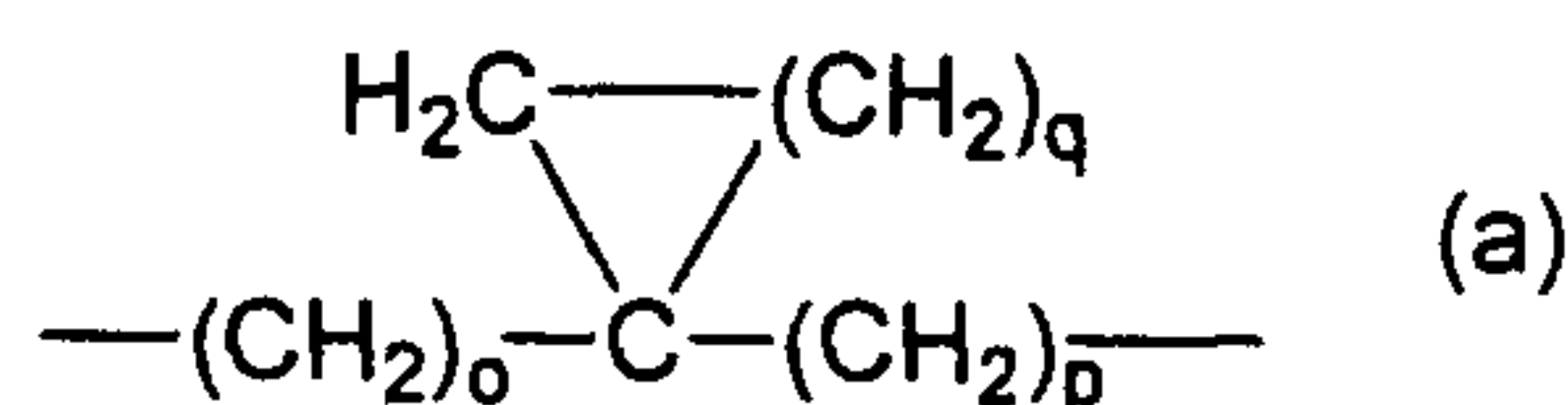
Summary of the invention

The present invention relates to a dibenzocycloheptene compound represented by the formula (I):



5 wherein R^1 represents a hydrogen atom, a halogen atom, a hydroxy group, a nitro group, a cyano group, a carbamoyl group, a formyl group, a carboxyl group, a 1H-tetrazol-5-yl group, a C_1 - C_4 alkyl group, a
 10 fluoro C_1 - C_4 alkyl group, a hydroxy C_1 - C_4 alkyl group, a C_2 - C_4 alkenyl group, a C_2 - C_4 alkynyl group, a C_1 - C_4 alkoxy group, a fluoro C_1 - C_4 alkoxy group, a C_1 - C_4 alkylthio group, a C_1 - C_4 alkylsulfinyl group or a C_1 - C_4 alkylsulfonyl group, R^2 represents a hydrogen
 15 atom, a halogen atom, a nitro group, a cyano group, a C_1 - C_4 alkyl group or a C_1 - C_4 alkoxy group, A represents a 5-membered or 6-membered heteroaromatic ring group containing 1 to 3 hetero atoms selected from the group consisting of a nitrogen atom, an oxygen
 20 atom and a sulfur atom or a fused heteroaromatic ring group in which the heteroaromatic ring group and a benzene ring are fused, the heteroaromatic ring group or fused heteroaromatic ring group may have a substituent(s) selected from a halogen atom, a nitro group, a cyano group, a C_1 - C_4 alkyl group, a
 25 fluoro C_1 - C_4 alkyl group, a C_1 - C_4 alkoxy group, a fluoro C_1 - C_4 alkoxy group, a C_1 - C_4 alkylthio group or a C_3 - C_4 alkylene group, B represents a formula:
 -CH=CH-, -CH₂O-, -CH₂CH₂-, -CH₂S-, -OCH₂- or -SCH₂-, X
 30 represents an oxygen atom, a sulfur atom, methylene group or a formula: =CH-, Y represents a C_1 - C_{10}

alkylene group which may have a substituent(s) selected from a halogen atom, C₁-C₄ alkyl group and C₁-C₄ alkoxy group, or a group represented by the formula (a):



wherein o and p each represent an integer of 0 to 2, and q represents an integer of 1 to 4, Z represents a carboxyl group which may be protected; 1H-tetrazol-5-yl group; the formula -NH-SO₂-R³; or the formula -CO-NH-SO₂-R³

wherein R³ represents a C₁-C₄ alkyl group, a fluoro C₁-C₄ alkyl group or a phenyl group which may have a halogen atom, a C₁-C₄ alkyl group, a fluoro C₁-C₄ alkyl group, a C₁-C₄ alkoxy group, a fluoro C₁-C₄ alkoxy group, a nitro group or a cyano group as a substituent(s),

m is an integer of 1 to 4, when m is 2 or more, a plural number of R¹ may be different from each other, n is an integer of 1 to 3, and when n is 2 or more, a plural number of R² may be different from each other,

----- represents a single bond or a double bond, or a pharmaceutically acceptable salt thereof.

Best mode for carrying out the invention

In the compound represented by the above-mentioned formula (I), as the halogen atom of R¹, there may be mentioned, for example, a fluorine atom, a chlorine atom, bromine atom or iodine atom, preferably a fluorine atom, a chlorine atom or bromine atom, more preferably a fluorine atom or a chlorine atom.

As the C₁-C₄ alkyl group of R¹, there may be mentioned, for example, a straight or branched C₁-C₄ alkyl group such as methyl, ethyl, propyl, isopropyl, butyl,

isobutyl, sec-butyl or t-butyl group, preferably methyl, ethyl, propyl or isopropyl group, more preferably methyl or ethyl group, particularly preferably methyl group.

As the fluoro C₁-C₄ alkyl group of R¹, there may be mentioned, for example, a straight or branched C₁-C₄ alkyl group substituted by 1 to 3 fluorine atoms such as fluoromethyl, difluoromethyl, trifluoromethyl, 2-fluoroethyl, 2,2,2-trifluoroethyl, 2-fluoropropyl, 3-fluoropropyl or 4-fluorobutyl group, preferably fluoromethyl, difluoromethyl, trifluoromethyl or 2-fluoroethyl group, more preferably fluoromethyl, difluoromethyl or trifluoromethyl group, particularly preferably difluoromethyl or trifluoromethyl group.

As the hydroxy C₁-C₄ alkyl group of R¹, there may be mentioned, for example, a straight or branched C₁-C₄ alkyl group substituted by hydroxy group such as a hydroxymethyl, 1-hydroxyethyl, 2-hydroxyethyl, 1-hydroxy-1-methylethyl, 1-hydroxypropyl, 2-hydroxypropyl, 3-hydroxypropyl, 1-hydroxybutyl or 4-hydroxybutyl group, preferably hydroxymethyl, 1-hydroxyethyl, 2-hydroxyethyl, 1-hydroxy-1-methylethyl, 1-hydroxypropyl or 2-hydroxypropyl group, more preferably hydroxymethyl, 1-hydroxyethyl, 1-hydroxy-1-methylethyl or 1-hydroxypropyl group, particularly preferably hydroxymethyl or 1-hydroxy-1-methylethyl group.

As the C₂-C₄ alkenyl group of R¹, there may be mentioned, for example, a straight or branched C₂-C₄ alkenyl group such as vinyl, 1-propenyl, allyl, isopropenyl, 1-butenyl, 2-butenyl, 3-butenyl, 1-methyl-1-propenyl, 2-methyl-1-propenyl or 2-methyl-2-propenyl group, preferably vinyl, 1-propenyl, allyl, 1-butenyl, 2-butenyl or 2-methyl-1-propenyl group, more preferably vinyl, 1-propenyl or allyl group, particularly preferably vinyl group.

As the C₂-C₄ alkynyl group of R¹, there may be mentioned, for example, a straight C₂-C₄ alkynyl group such as ethynyl, 1-propynyl, propargyl, 1-butyne, 2-butyne or 3-butyne group, preferably ethynyl, 1-propynyl or 1-butyne

group, more preferably ethynyl or 1-propynyl group, particularly preferably ethynyl group.

As the C₁-C₄ alkoxy group of R¹, there may be mentioned, for example, a straight or branched C₁-C₄ alkoxy group such as methoxy, ethoxy, propoxy, isopropoxy, butoxy, isobutoxy, sec-butoxy or t-butoxy group, preferably methoxy, ethoxy, propoxy or isopropoxy group, more preferably methoxy or ethoxy group, particularly preferably methoxy group.

As the fluoro C₁-C₄ alkoxy group of R¹, there may be mentioned, for example, a straight or branched C₁-C₄ alkoxy group substituted by 1 to 3 fluorine atoms such as fluoromethoxy, difluoromethoxy, trifluoromethoxy, 2-fluoroethoxy, 2, 2, 2-trifluoroethoxy, 2-fluoropropoxy, 3-fluoropropoxy or 4-fluorobutoxy group, preferably fluoromethoxy, difluoromethoxy, trifluoromethoxy or 2-fluoroethoxy group, more preferably fluoromethoxy, difluoromethoxy or trifluoromethoxy group, particularly preferably difluoromethoxy or trifluoromethoxy group.

As the C₁-C₄ alkylthio group of R¹, there may be mentioned, for example, a straight or branched C₁-C₄ alkylthio group such as methylthio, ethylthio, propylthio, isopropylthio, butylthio, isobutylthio, sec-butylthio or t-butylthio group, preferably methylthio, ethylthio, propylthio or isopropylthio group, more preferably methylthio or ethylthio group, particularly preferably methylthio group.

As the C₁-C₄ alkylsulfinyl group of R¹, there may be mentioned, for example, a straight or branched C₁-C₄ alkylsulfinyl group such as methylsulfinyl, ethylsulfinyl, propylsulfinyl, isopropylsulfinyl, butylsulfinyl, isobutylsulfinyl, sec-butylsulfinyl or t-butylsulfinyl group, preferably methylsulfinyl, ethylsulfinyl, propylsulfinyl or isopropylsulfinyl group, more preferably methylsulfinyl or ethylsulfinyl group, particularly preferably methylsulfinyl group.

As the C₁-C₄ alkylsulfonyl group of R¹, there may be

mentioned, for example, a straight or branched C₁-C₄ alkyl-sulfonyl group such as methylsulfonyl, ethylsulfonyl, propylsulfonyl, isopropylsulfonyl, butylsulfonyl, isobutylsulfonyl, sec-butylsulfonyl or t-butylsulfonyl group, preferably methylsulfonyl, ethylsulfonyl, propylsulfonyl or isopropylsulfonyl group, more preferably methylsulfonyl or ethylsulfonyl group, particularly preferably methylsulfonyl group.

In particular, as R¹ in the formula (I), there may be preferably mentioned a hydrogen atom, a fluorine atom, a chlorine atom, bromine atom, a hydroxy group, a nitro group, a cyano group, a carbamoyl group, a formyl group, a carboxyl group, 1H-tetrazol-5-yl group, methyl group, ethyl group, propyl group, isopropyl group, fluoromethyl group, difluoromethyl group, trifluoromethyl group, 2-fluoroethyl group, hydroxymethyl group, 1-hydroxyethyl group, 2-hydroxyethyl group, 1-hydroxy-1-methylethyl group, 1-hydroxypropyl group, 2-hydroxypropyl group, vinyl group, 1-propenyl group, allyl group, 1-butenyl group, 2-butenyl group, 2-methyl-1-propenyl group, ethynyl group, 1-propynyl group, 1-butylnyl group, methoxy group, ethoxy group, propoxy group, isopropoxy group, fluoromethoxy group, difluoromethoxy group, trifluoromethoxy group, 2-fluoroethoxy group, methylthio group, ethylthio group, propylthio group, isopropylthio group, methylsulfinyl group, ethylsulfinyl group, propylsulfinyl group, isopropylsulfinyl group, methylsulfonyl group, ethylsulfonyl group, propylsulfonyl group or isopropylsulfonyl group,

more preferably a hydrogen atom, a fluorine atom, a chlorine atom, a hydroxy group, a nitro group, a cyano group, a carbamoyl group, a formyl group, 1H-tetrazol-5-yl group, methyl group, ethyl group, fluoromethyl group, difluoromethyl group, trifluoromethyl group, hydroxymethyl group, 1-hydroxyethyl group, 1-hydroxy-1-methylethyl group, 1-hydroxypropyl group, vinyl group, 1-propenyl group, allyl group, ethynyl group, 1-propynyl group, 1-butylnyl group,

methoxy group, ethoxy group, fluoromethoxy group, difluoromethoxy group, trifluoromethoxy group, methylthio group, ethylthio group, methylsulfinyl group, ethylsulfinyl group, methylsulfonyl group or ethylsulfonyl group, further more preferably hydrogen atom, fluorine atom, chlorine atom, nitro group, cyano group, formyl group, 1H-tetrazol-5-yl group, methyl group, difluoromethyl group, trifluoromethyl group, hydroxymethyl group, 1-hydroxy-1-methylethyl group, vinyl group, ethynyl group, methoxy group, difluoromethoxy group, trifluoromethoxy group, methylthio group, methylsulfinyl group or methylsulfonyl group, particularly preferably hydrogen atom, fluorine atom, chlorine atom, cyano group, trifluoromethyl group, hydroxymethyl group, 1-hydroxy-1-methylethyl group, ethynyl group, methoxy group, difluoromethoxy group, trifluoromethoxy group, methylsulfinyl group or methylsulfonyl group.

In the formula (I), a halogen atom, C₁-C₄ alkyl group and C₁-C₄ alkoxy group of R² have the same meanings as those of the above-mentioned R¹, and as R², there may be preferably mentioned hydrogen atom, fluorine atom, chlorine atom, bromine atom, nitro group, cyano group, methyl group, ethyl group, propyl group, isopropyl group, methoxy group, ethoxy group, propoxy group or isopropoxy group, more preferably hydrogen atom, fluorine atom, chlorine atom, nitro group, cyano group, methyl group, ethyl group, methoxy group or ethoxy group, further more preferably hydrogen atom, fluorine atom, chlorine atom, methyl group or methoxy group, particularly preferably hydrogen atom.

In the formula (I), "a 5-membered or 6-membered heteroaromatic ring group containing 1 to 3 hetero atom(s) selected from the group consisting of a nitrogen atom, an oxygen atom and a sulfur atom or a fused heteroaromatic ring group in which the heteroaromatic ring group and a benzene are fused" of A, there may be mentioned, for example, a 5-membered heteroaromatic ring group such as furan, thiophen, oxazol, thiazol, imidazol, pyrazol or

thiadiazol group; a 6-membered heteroaromatic ring group such as pyridin, pyrimidin, pyridazin or pyrazin group; or a fused heteroaromatic ring group such as benzofuran, benzothiophen, benzoxazol, benzothiazol, benzimidazol, quinolin, quinazolin or quinoxalin group, preferably oxazol, thiazol, imidazol, pyrazol, thiadiazol, pyridine, pyrimidin, pyridazin, pyrazin, benzoxazol, benzothiazol, benzimidazol, quinolin, quinazolin or quinoxalin group, more preferably thiazol, thiadiazol, pyridine, pyrimidin, benzoxazol, benzothiazol, quinolin or quinazolin group, particularly preferably pyridine, benzothiazol or quinolin group.

The above-mentioned heteroaromatic ring group or fused heteroaromatic ring group may have a substituent(s), and as the substituent(s), there may be mentioned, for example, a halogen atom with the same meaning as R^1 ; a C_1-C_4 alkyl group with the same meaning as R^1 ; a fluoro C_1-C_4 alkyl group with the same meaning as R^1 ; a C_1-C_4 alkoxy group with the same meaning as R^1 ; a fluoro C_1-C_4 alkoxy group with the same meaning as R^1 ; a C_1-C_4 alkylthio group with the same meaning as R^1 ; nitro group; cyano group; or a C_3-C_4 alkylene group such as trimethylene, tetramethylene group (said alkylene group bonds to an adjacent carbon atom on the heteroaromatic ring to form a 5-membered ring or 6-membered ring), preferably fluorine atom, chlorine atom, bromine atom, nitro, cyano, methyl, ethyl, propyl, isopropyl, butyl, isobutyl, t-butyl, fluoromethyl, difluoromethyl, trifluoromethyl, 2-fluoroethyl, methoxy, ethoxy, propoxy, isopropoxy, fluoromethoxy, difluoromethoxy, trifluoromethoxy, 2-fluoroethoxy, methylthio, ethylthio, propylthio, isopropylthio, trimethylene or tetramethylene group, more preferably a fluorine atom, a chlorine atom, nitro, cyano, methyl, ethyl, isopropyl, t-butyl, fluoromethyl, difluoromethyl, trifluoromethyl, methoxy, ethoxy, fluoromethoxy, difluoromethoxy, trifluoromethoxy, methylthio, ethylthio, trimethylene or tetramethylene group,

further more preferably fluorine atom, chlorine atom, nitro, cyano, methyl, isopropyl, t-butyl, difluoromethyl, trifluoromethyl, methoxy, difluoromethoxy, trifluoromethoxy, methylthio or tetramethylene group, particularly preferably fluorine atom, chlorine atom, trifluoromethyl or tetramethylene group.

A number of the substituent(s) on the heteroaromatic ring group or fused heteroaromatic ring group is 1 to 4, preferably 1 to 2.

As A in the formula (I), it is specifically mentioned, preferably 2-oxazolyl, 2-thiazolyl, 2- or 4-imidazolyl, 3-pyrazolyl, 1,3,4-thiadiazol-2-yl, 2-pyridyl, 2- or 4-pyrimidinyl, 3-pyridazinyl, 2-pyrazinyl, 2-benzoxazolyl, 2-benzothiazolyl, 2-benzoimidazolyl, quinolin-2-yl, quinazolin-2-yl, quinoxalin-2-yl, 4-methyl-2-thiazolyl, 4-ethyl-2-thiazolyl, 4-isopropyl-2-thiazolyl, 4-t-butyl-2-thiazolyl, 4-trifluoromethyl-2-thiazolyl, 5-methyl-1,3,4-thiadiazol-2-yl, 5-ethyl-1,3,4-thiadiazol-2-yl, 5-isopropyl-1,3,4-thiadiazol-2-yl, 5-t-butyl-1,3,4-thiadiazol-2-yl, 5-trifluoromethyl-1,3,4-thiadiazol-2-yl, 5,6-difluoro-2-pyridyl, 5,6-dichloro-2-pyridyl, 5,6-dimethyl-2-pyridyl, 5,6-diethyl-2-pyridyl, 6-trifluoromethyl-2-pyridyl, 6-methylthio-2-pyridyl, 5H-6,7-dihydrocyclopenta[b]pyridin-2-yl, 5,6,7,8-tetrahydroquinolin-2-yl, 5,6-difluoro-2-pyrimidinyl, 5,6-dichloro-2-pyrimidinyl, 5,6-dimethyl-2-pyrimidinyl, 6-trifluoromethyl-2-pyrimidinyl, 5H-6,7-dihydrocyclopenta[d]pyrimidin-2-yl, 5,6,7,8-tetrahydroquinazolin-2-yl, 6-fluoro-2-benzoxazolyl, 5-fluoro-2-benzoxazolyl, 5,6-difluoro-2-benzoxazolyl, 6-chloro-2-benzoxazolyl, 5-chloro-2-benzoxazolyl, 5,6-dichloro-2-benzoxazolyl, 5-chloro-6-fluoro-2-benzoxazolyl, 5-methyl-2-benzoxazolyl, 5-cyano-2-benzoxazolyl, 5-trifluoromethyl-2-benzoxazolyl, 5-methylthio-2-benzoxazolyl, 6-fluoro-2-benzothiazolyl, 5-fluoro-2-benzothiazolyl, 5,6-difluoro-2-benzothiazolyl, 6-chloro-2-benzothiazolyl, 5-chloro-2-benzothiazolyl, 5,6-dichloro-2-benzothiazolyl, 5-chloro-6-

fluoro-2-benzothiazolyl, 5-methyl-2-benzothiazolyl, 5-cyano-2-benzothiazolyl, 5-trifluoromethyl-2-benzothiazolyl, 5-methylthio-2-benzothiazolyl, 5-fluoroquinolin-2-yl, 6-fluoroquinolin-2-yl, 7-fluoroquinolin-2-yl, 5-chloro-quinolin-2-yl, 6-chloroquinolin-2-yl, 7-chloroquinolin-2-yl, 7-methylquinolin-2-yl, 7-trifluoromethylquinolin-2-yl, 7-methoxyquinolin-2-yl, 7-difluoromethoxyquinolin-2-yl, 7-trifluoromethoxyquinolin-2-yl, 5,7-difluoroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 5,7-dichloroquinolin-2-yl, 6,7-dichloroquinolin-2-yl, 5-chloro-7-fluoroquinolin-2-yl, 6-chloro-7-fluoroquinolin-2-yl, 7-chloro-5-fluoroquinolin-2-yl, 7-chloro-6-fluoroquinolin-2-yl, 7-chloro-6-cyanoquinolin-2-yl, 7-cyano-6-fluoroquinolin-2-yl, 6-fluoro-7-trifluoromethylquinolin-2-yl, 5,6,7-trifluoroquinolin-2-yl, 5-fluoroquinazolin-2-yl, 6-fluoroquinazolin-2-yl, 7-fluoroquinazolin-2-yl, 5-chloroquinazolin-2-yl, 6-chloroquinazolin-2-yl, 7-chloroquinazolin-2-yl, 7-methylquinazolin-2-yl, 7-trifluoromethylquinazolin-2-yl, 7-methoxyquinazolin-2-yl, 7-difluoromethoxyquinazolin-2-yl, 7-trifluoromethoxyquinazolin-2-yl, 5,7-difluoroquinazolin-2-yl, 6,7-difluoroquinazolin-2-yl, 5,7-dichloroquinazolin-2-yl, 6,7-dichloroquinazolin-2-yl, 5-chloro-7-fluoroquinazolin-2-yl, 6-chloro-7-fluoroquinazolin-2-yl, 7-chloro-5-fluoroquinazolin-2-yl, 7-chloro-6-fluoroquinazolin-2-yl, 7-chloro-6-cyanoquinazolin-2-yl, 7-cyano-6-fluoroquinazolin-2-yl, 6-fluoro-7-trifluoromethylquinazolin-2-yl or 5,6,7-trifluoroquinazolin-2-yl group,

more preferably 2-thiazolyl, 1,3,4-thiadiazol-2-yl, 2-pyridyl, 2-pyrimidinyl, 2-benzoxazolyl, 2-benzothiazolyl, quinolin-2-yl, quinazolin-2-yl, 4-methyl-2-thiazolyl, 4-isopropyl-2-thiazolyl, 4-t-butyl-2-thiazolyl, 4-trifluoromethyl-2-thiazolyl, 5-methyl-1,3,4-thiadiazol-2-yl, 5-isopropyl-1,3,4-thiadiazol-2-yl, 5-t-butyl-1,3,4-thiadiazol-2-yl, 5-trifluoromethyl-1,3,4-thiadiazol-2-yl, 5,6-difluoro-2-pyridyl, 5,6-dichloro-2-pyridyl, 5,6-dimethyl-2-pyridyl, 5H-6,7-dihydrocyclopenta[b]pyridin-2-yl, 5,6,7,8-

tetrahydroquinolin-2-yl, 5,6-difluoro-2-pyrimidinyl, 5,6-dichloro-2-pyrimidinyl, 5,6-dimethyl-2-pyrimidinyl, 6-trifluoromethyl-2-pyrimidinyl, 5H-6,7-dihydrocyclopenta-[d]pyrimidin-2-yl, 5,6,7,8-tetrahydroquinazolin-2-yl, 6-fluoro-2-benzoxazolyl, 5-fluoro-2-benzoxazolyl, 5,6-difluoro-2-benzoxazolyl, 6-chloro-2-benzoxazolyl, 5-chloro-2-benzoxazolyl, 5,6-dichloro-2-benzoxazolyl, 5-chloro-6-fluoro-2-benzoxazolyl, 5-methyl-2-benzoxazolyl, 5-cyano-2-benzoxazolyl, 5-trifluoromethyl-2-benzoxazolyl, 5-methylthio-2-benzoxazolyl, 6-fluoro-2-benzothiazolyl, 5-fluoro-2-benzothiazolyl, 5,6-difluoro-2-benzothiazolyl, 6-chloro-2-benzothiazolyl, 5-chloro-2-benzothiazolyl, 5,6-dichloro-2-benzothiazolyl, 5-chloro-6-fluoro-2-benzothiazolyl, 5-methyl-2-benzothiazolyl, 5-cyano-2-benzothiazolyl, 5-trifluoromethyl-2-benzothiazolyl, 5-methylthio-2-benzothiazolyl, 5-fluoroquinolin-2-yl, 6-fluoroquinolin-2-yl, 7-fluoroquinolin-2-yl, 5-chloroquinolin-2-yl, 6-chloroquinolin-2-yl, 7-chloroquinolin-2-yl, 7-methylquinolin-2-yl, 7-trifluoromethylquinolin-2-yl, 7-methoxyquinolin-2-yl, 7-difluoromethoxyquinolin-2-yl, 7-trifluoromethoxyquinolin-2-yl, 5,7-difluoroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 5,7-dichloroquinolin-2-yl, 6,7-dichloroquinolin-2-yl, 5-chloro-7-fluoroquinolin-2-yl, 6-chloro-7-fluoroquinolin-2-yl, 7-chloro-5-fluoroquinolin-2-yl, 7-chloro-6-fluoroquinolin-2-yl, 7-chloro-6-cyanoquinolin-2-yl, 7-cyano-6-fluoroquinolin-2-yl, 6-fluoro-7-trifluoromethylquinolin-2-yl, 5,6,7-trifluoroquinolin-2-yl, 5-fluoroquinazolin-2-yl, 6-fluoroquinazolin-2-yl, 7-fluoroquinazolin-2-yl, 5-chloroquinazolin-2-yl, 6-chloroquinazolin-2-yl, 7-chloroquinazolin-2-yl, 7-methylquinazolin-2-yl, 7-trifluoromethylquinazolin-2-yl, 7-methoxyquinazolin-2-yl, 7-difluoromethoxyquinazolin-2-yl, 7-trifluoromethoxyquinazolin-2-yl, 5,7-difluoroquinazolin-2-yl, 6,7-difluoroquinazolin-2-yl, 5,7-dichloroquinazolin-2-yl, 6,7-dichloroquinazolin-2-yl, 5-chloro-7-fluoroquinazolin-2-yl, 6-chloro-7-fluoroquinazolin-2-yl, 7-chloro-5-fluoroquinazolin-2-yl, 7-chloro-6-

fluoroquinazolin-2-yl, 7-chloro-6-cyanoquinazolin-2-yl, 7-cyano-6-fluoroquinazolin-2-yl, 6-fluoro-7-trifluoromethylquinazolin-2-yl or 5,6,7-trifluoroquinazolin-2-yl group,
 further more preferably 2-pyridyl, 2-benzothiazolyl,
 5 quinolin-2-yl, 5,6-difluoro-2-pyridyl, 5,6-dichloro-2-pyridyl, 5,6-dimethyl-2-pyridyl, 5,6,7,8-tetrahydroquinolin-2-yl, 6-fluoro-2-benzothiazolyl, 5-fluoro-2-benzothiazolyl, 5,6-difluoro-2-benzothiazolyl, 6-chloro-2-benzothiazolyl, 5-chloro-2-benzothiazolyl, 5,6-dichloro-2-benzothiazolyl, 5-chloro-6-fluoro-2-benzothiazolyl, 5-methyl-2-benzothiazolyl, 5-cyano-2-benzothiazolyl, 5-trifluoromethyl-2-benzothiazolyl, 5-methylthio-2-benzothiazolyl, 5-fluoroquinolin-2-yl, 6-fluoroquinolin-2-yl, 7-fluoroquinolin-2-yl, 5-chloroquinolin-2-yl, 6-chloroquinolin-2-yl, 7-chloroquinolin-2-yl, 7-methylquinolin-2-yl, 7-trifluoromethylquinolin-2-yl, 7-methoxyquinolin-2-yl, 7-difluoromethoxyquinolin-2-yl, 7-trifluoromethoxyquinolin-2-yl, 5,7-difluoroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 5,7-dichloroquinolin-2-yl, 6,7-dichloroquinolin-2-yl, 5-chloro-7-fluoroquinolin-2-yl, 6-chloro-7-fluoroquinolin-2-yl, 7-chloro-5-fluoroquinolin-2-yl, 7-chloro-6-fluoroquinolin-2-yl, 7-chloro-6-cyanoquinolin-2-yl, 7-cyano-6-fluoroquinolin-2-yl, 6-fluoro-7-trifluoromethylquinolin-2-yl or 5,6,7-trifluoroquinolin-2-yl group,
 10 particularly preferably 5,6,7,8-tetrahydroquinolin-2-yl, 7-fluoroquinolin-2-yl, 7-chloroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 6,7-dichloroquinolin-2-yl or 7-chloro-6-fluoroquinolin-2-yl.
 25

In the above-mentioned formula (I), B is, for
 30 example, a formula: $-\text{CH}=\text{CH}-$, a formula: $-\text{CH}_2\text{O}-$, a formula: $-\text{CH}_2\text{CH}_2-$, a formula: $-\text{CH}_2\text{S}-$, a formula: $-\text{OCH}_2-$ or a formula: $-\text{SCH}_2-$, preferably a formula: $-\text{CH}=\text{CH}-$, a formula: $-\text{OCH}_2-$ or a formula: $-\text{CH}_2\text{O}-$.

In the above-mentioned formula (I), X is an oxygen
 35 atom, a sulfur atom, a methylene group or a formula: $=\text{CH}-$, preferably an oxygen atom or a sulfur atom.

As the C₁-C₁₀ alkylene group of Y in the above-mentioned formula (I), there may be mentioned, for example, a straight C₁-C₁₀ alkylene group such as methylene, ethylene, trimethylene, tetramethylene, pentamethylene, hexamethylene, heptamethylene, octamethylene, nonamethylene or decamethylene group, preferably a C₁-C₆ alkylene group, more preferably a C₁-C₄ alkylene group,

particularly preferably methylene, ethylene or trimethylene group.

Said alkylene group may have a substituent(s), and a halogen atom, a C₁-C₄ alkyl group and a C₁-C₄ alkoxy group as said substituent(s) thereof have the same meanings as those mentioned in the above-mentioned R¹, respectively.

As the substituent(s) for the alkylene group of Y, there may be preferably mentioned fluorine atom, chlorine atom, methyl, ethyl, propyl, methoxy, ethoxy or propoxy group, more preferably fluorine atom, methyl, ethyl or methoxy group, particularly preferably fluorine atom or methyl group.

As the group represented by the formula (a) of Y, there may be preferably mentioned a group wherein o=0, p=0 and q=1 (hereinafter referred to as (a-1) group), a group wherein o=0, p=1 and q=1 (hereinafter referred to as (a-2) group), a group wherein o=0, p=1 and q=2 (hereinafter referred to as (a-3) group), a group wherein o=1, p=0 and q=1 (hereinafter referred to as (a-4) group), a group wherein o=1, p=1 and q=1 (hereinafter referred to as (a-5) group), a group wherein o=1, p=1 and q=2 (hereinafter referred to as (a-6) group) or a group wherein o=1, p=1 and q=3 (hereinafter referred to as (a-7) group), more preferably a (a-4) group, a (a-5) group or a (a-6) group, particularly more preferably (a-5) group.

As the preferred group of Y in the formula (I), there may be specifically mentioned methylene, ethylene, trimethylene, tetramethylene, pentamethylene, fluoromethylene, difluoromethylene, 1-fluoroethylene, 2-fluoro-

ethylene, 1,1-difluoroethylene, 2,2-difluoroethylene, ethylidene, 1-methylethylene, 2-methylethylene, 1,1-dimethylethylene, 2,2-dimethylethylene, 1-ethylethylene, 2-ethylethylene, 1-methoxyethylene, 2-methoxyethylene, 1-fluorotrimethylene, 2-fluorotrimethylene, 3-fluorotrimethylene, 1,1-difluorotrimethylene, 2,2-difluorotrimethylene, 3,3-difluorotrimethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethyltrimethylene, 2,2-dimethyltrimethylene, 3,3-dimethyltrimethylene, 2,2-diethyltrimethylene, 2-methoxytrimethylene, 3-methoxytrimethylene, 2,2-dimethoxytrimethylene, 3,3-dimethoxytrimethylene, (a-1) group, (a-2) group, (a-3) group, (a-4) group, (a-5) group or (a-6) group, more preferably methylene, ethylene, trimethylene, fluoromethylene, difluoromethylene, 1-fluoroethylene, 2-fluoroethylene, 1,1-difluoroethylene, 2,2-difluoroethylene, ethylidene, 1-methylethylene, 2-methylethylene, 1-fluorotrimethylene, 2-fluorotrimethylene, 3-fluorotrimethylene, 1,1-difluorotrimethylene, 2,2-difluorotrimethylene, 3,3-difluorotrimethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethyltrimethylene, 2,2-dimethyltrimethylene, 3,3-dimethyltrimethylene, (a-4) group, (a-5) group or (a-6) group, further more preferably methylene, ethylene, trimethylene, difluoromethylene, 1-fluoroethylene, 2-fluoroethylene, 1,1-difluoroethylene, 2,2-difluoroethylene, ethylidene, 1-methylethylene, 2-methylethylene, 2,2-difluorotrimethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethyltrimethylene, 2,2-dimethyltrimethylene, 3,3-dimethyltrimethylene or (a-5) group, particularly preferably methylene, ethylene, trimethylene, ethylidene, 1-methylethylene, 2-methylethylene or (a-5) group.

In a group of the formula: $\text{-NH-SO}_2\text{-R}^3$ or the formula: $\text{-CO-NH-SO}_2\text{-R}^3$ group shown by Z, $\text{C}_1\text{-C}_4$ alkyl group of R^3 ; a fluoro $\text{C}_1\text{-C}_4$ alkyl group of R^3 ; or a halogen atom, $\text{C}_1\text{-C}_4$ alkyl group, fluoro $\text{C}_1\text{-C}_4$ alkyl group, $\text{C}_1\text{-C}_4$ alkoxy group or fluoro $\text{C}_1\text{-C}_4$ alkoxy group which is a substituent(s) on a

phenyl group of R^3 have the same meanings as those mentioned in the above-mentioned R^1 , respectively.

As R^3 , there may be preferably mentioned methyl, ethyl, propyl, fluoromethyl, difluoromethyl, trifluoromethyl, 2-fluoroethyl, 2, 2, 2-trifluoroethyl, phenyl, (o-, m- or p-)fluorophenyl, (o-, m- or p-)chlorophenyl, (o-, m- or p-)methylphenyl, (o-, m- or p-)ethylphenyl, (o-, m- or p-)(trifluoromethyl)phenyl, (o-, m- or p-)methoxyphenyl, (o-, m- or p-)ethoxyphenyl, (o-, m- or p-)(difluoromethoxy)phenyl, (o-, m- or p-)(trifluoromethoxy)phenyl, (o-, m- or p-)nitrophenyl or (o-, m- or p-)cyanophenyl group, more preferably methyl, ethyl, trifluoromethyl, 2-fluoroethyl, 2,2,2-trifluoroethyl, phenyl, (o- or p-)fluorophenyl, (o- or p-)chlorophenyl, (o- or p-)methylphenyl, (o- or p-)(trifluoromethyl)phenyl, (o- or p-)methoxyphenyl, (o- or p-)(difluoromethoxy)phenyl, (o- or p-)(trifluoromethoxy)phenyl, (o- or p-)nitrophenyl or (o- or p-)cyanophenyl group, further more preferably methyl, ethyl, trifluoromethyl, phenyl, p-fluorophenyl, p-chlorophenyl, (o- or p-)methylphenyl, p-(trifluoromethyl)phenyl, (o- or p-)methoxyphenyl, p-(difluoromethoxy)phenyl, p-(trifluoromethoxy)phenyl, p-nitrophenyl or p-cyanophenyl group, particularly preferably methyl, trifluoromethyl, phenyl, o-methylphenyl or p-methylphenyl group.

As a preferred group of Z in the formula (I), there may be specifically mentioned carboxy, 1H-tetrazol-5-yl, methanesulfonylamino, ethanesulfonylamino, trifluoromethanesulfonylamino, phenylsulfonylamino, p-fluorophenylsulfonylamino, p-chlorophenylsulfonylamino, o-methylphenylsulfonylamino, p-methylphenylsulfonylamino, p-trifluoromethylphenylsulfonylamino, o-methoxyphenylsulfonylamino, p-methoxyphenylsulfonylamino, p-difluoromethoxyphenylsulfonylamino, p-trifluoromethoxyphenylsulfonylamino, p-nitrophenylsulfonylamino, p-cyanophenylsulfonylamino, methanesulfonylaminocarbonyl, ethanesulfonylaminocarbonyl, trifluoromethanesulfonylaminocarbonyl, phenylsulfonyl-

aminocarbonyl, p-fluorophenylsulfonylamino-
 carbonyl, p-chlorophenylsulfonylamino-
 carbonyl, o-methylphenylsulfonyl-
 aminocarbonyl, p-methylphenylsulfonylamino-
 carbonyl, p-tri-
 fluoromethylphenylsulfonylamino-
 carbonyl, o-methoxyphenyl-
 5 sulfonylamino-
 carbonyl, p-methoxyphenylsulfonylamino-
 carbonyl, p-difluoromethoxyphenylsulfonylamino-
 carbonyl, p-tri-
 fluoromethoxyphenylsulfonylamino-
 carbonyl, p-nitrophenyl-
 sulfonylamino-
 carbonyl or p-cyanophenylsulfonylamino-
 carbonyl group,

10 more preferably carboxy, 1H-tetrazol-5-yl, methane-
 sulfonylamino, trifluoromethanesulfonylamino, phenylsulfonyl-
 amino, o-methylphenylsulfonylamino, p-methylphenylsulfonyl-
 amino, methanesulfonylamino-
 carbonyl, trifluoromethane-
 sulfonylamino-
 carbonyl, phenylsulfonylamino-
 carbonyl, o-
 15 methylphenylsulfonylamino-
 carbonyl or p-methylphenylsulfonyl-
 amino-
 carbonyl group, further more preferably carboxy,
 methanesulfonylamino, trifluoromethanesulfonylamino,
 methanesulfonylamino-
 carbonyl or trifluoromethanesulfonyl-
 amino-
 carbonyl group, particularly preferably carboxyl
 20 group.

Incidentally, when Z is a carboxyl group, the
 carboxyl group may be protected by a protective group. As
 the protective group, it is not specifically limited so
 long as it can be easily deprotected in a living body to be
 25 changed to a carboxyl group, and there may be mentioned,
 for example, a C₁-C₄ alkyl group having the same meanings as
 those defined in R¹; a C₇-C₁₀ aralkyl group such as benzyl,
 phenylethyl or phenylpropyl group; a C₁-C₄ alkyl group
 substituted by a C₂-C₅ alkanoyloxy group, such as acetoxy-
 30 methyl, 1-acetoxyethyl, 1-acetoxypropyl, 1-acetoxybutyl,
 propanoyloxymethyl, 1-propanoyloxyethyl, butanoyloxymethyl,
 1-butanoyloxymethyl, pivaloyloxymethyl, 1-pivaloyloxyethyl,
 1-pivaloyloxypropyl or 1-pivaloyloxybutyl group; a C₁-C₄
 alkyl group substituted by a (C₁-C₄ alkoxy)carbonyloxy
 35 group, such as methoxycarbonyloxymethyl, 1-(methoxycarbonyl-
 oxy)ethyl, ethoxycarbonyloxymethyl, 1-(ethoxycarbonyl-

oxy)ethyl, propoxycarbonyloxymethyl, oxymethylcarbonyloxymethyl, 1-(isopropoxycarbonyloxy)ethyl, butoxycarbonyloxymethyl, 1-(butoxycarbonyloxy)ethyl, t-butoxycarbonyloxymethyl or 1-(t-butoxycarbonyloxy)ethyl group; a C₁-C₄ alkyl group substituted by a N,N-di(C₁-C₄ alkyl)amino-carbonyl group such as N,N-dimethylaminocarbonylmethyl, 2-(N,N-dimethylaminocarbonyl)ethyl or N,N-diethylaminocarbonylmethyl group; a C₁-C₄ alkyl group substituted by a N,N-di(C₁-C₄ alkyl)amino group or a 5- to 6-membered cyclic amino group which may contain an oxygen atom such as 2-(N,N-dimethylamino)ethyl, 2-(N,N-diethylamino)ethyl, 3-(N,N-dimethylamino)propyl, 2-piperidinoethyl, 2-(4-methyl)-piperidinoethyl, 3-piperidinopropyl or 2-morpholinoethyl group; or (5-methyl-2-oxo-1,3-dioxolen-4-yl)methyl group, and the like.

As a protective group for the carboxyl group, it is preferably a C₁-C₄ alkyl group; benzyl group; a C₁-C₂ alkyl group substituted by a C₂-C₅ alkanoyloxy group; a C₁-C₂ alkyl group substituted by a (C₁-C₄ alkoxy)carbonyloxy group; or (5-methyl-2-oxo-1,3-dioxolen-4-yl)methyl group, more preferably methyl, ethyl, propyl, isopropyl, butyl, isobutyl, acetoxymethyl, 1-acetoxyethyl, pivaloyloxymethyl or 1-pivaloyloxyethyl group.

In the above-mentioned formula (I), m is an integer of 1 to 4, m is preferably 1, 2 or 3, and particularly preferably 1 or 2. When m is 2 or more, R¹s may be different from each other.

In the above-mentioned formula (I), n is an integer of 1 to 3, preferably n is 1 or 2, particularly preferably 1. When n is 2 or more, R²s may be different from each other.

In the Compound (I) of the present invention, there exist an optical isomer(s) (including diastereomer) due to an asymmetric carbon atom(s) in the molecule, or there exist a case in which a geometric isomer due to a double bond exists, and these respective isomers are also included

in the present invention.

Also, the Compound (I) of the present invention can be converted into a pharmaceutically acceptable salt, if necessary. Such a pharmaceutically acceptable salt may be mentioned an acid addition salt of a mineral acid such as hydrochloride, hydrobromide, hydroiodide, sulfate or phosphate; an acid addition salt of an organic acid such as trifluoroacetate, methanesulfonate, ethanesulfonate, benzenesulfonate, p-toluenesulfonate, oxalate, maleate, fumarate, tartarate or citrate; a metal salt of a carboxylic acid such as a sodium salt, a potassium salt, a calcium salt, a magnesium salt, a manganese salt, an iron salt or an aluminum salt; or a salt with an organic base such as an ammonium salt, a triethylamine salt, a guanidine salt, a hydrazine salt, a quinine salt or a cinchonine salt, and the like.

Incidentally, the Compound (I) of the present invention can also exist as a hydrate.

In the dibenzocycloheptene compound having the above-mentioned formula (I) of the present invention, it is preferably

(1). a dibenzocycloheptene compound wherein R^1 of the compound represented by the formula (I) is selected from the group consisting of hydrogen atom, fluorine atom, chlorine atom, bromine atom, hydroxy group, nitro group, a cyano group, carbamoyl group, formyl group, carboxyl group, 1H-tetrazol-5-yl group, methyl group, ethyl group, propyl group, isopropyl group, fluoromethyl group, difluoromethyl group, trifluoromethyl group, 2-fluoroethyl group, hydroxymethyl group, 1-hydroxyethyl group, 2-hydroxyethyl group, 1-hydroxy-1-methylethyl group, 1-hydroxypropyl group, 2-hydroxypropyl group, vinyl group, 1-propenyl group, allyl group, 1-butenyl group, 2-butenyl group, 2-methyl-1-propenyl group, ethynyl group, 1-propynyl group, 1-butylnyl group, methoxy group, ethoxy group, propoxy group, isopropoxy group, fluoromethoxy group,

- difluoromethoxy group, trifluoromethoxy group, 2-fluoroethoxy group, methylthio group, ethylthio group, propylthio group, isopropylthio group, methylsulfinyl group, ethylsulfinyl group, propylsulfinyl group, isopropylsulfinyl group, methylsulfonyl group, ethylsulfonyl group, propylsulfonyl group and isopropylsulfonyl group,
- (2). a dibenzocycloheptene compound wherein R^1 of the compound represented by the formula (I) is selected from the group consisting of hydrogen atom, fluorine atom, chlorine atom, hydroxy group, nitro group, cyano group, carbamoyl group, formyl group, 1H-tetrazol-5-yl group, methyl group, ethyl group, fluoromethyl group, difluoromethyl group, trifluoromethyl group, hydroxymethyl group, 1-hydroxyethyl group, 1-hydroxy-1-methylethyl group, 1-hydroxypropyl group, vinyl group, 1-propenyl group, allyl group, ethynyl group, 1-propynyl group, 1-butynyl group, methoxy group, ethoxy group, fluoromethoxy group, difluoromethoxy group, trifluoromethoxy group, methylthio group, ethylthio group, methylsulfinyl group, ethylsulfinyl group, methylsulfonyl group and ethylsulfonyl group,
- (3). a dibenzocycloheptene compound wherein R^1 of the compound represented by the formula (I) is selected from the group consisting of a hydrogen atom, a fluorine atom, a chlorine atom, a nitro group, a cyano group, a formyl group, 1H-tetrazol-5-yl group, methyl group, difluoromethyl group, trifluoromethyl group, hydroxymethyl group, 1-hydroxy-1-methylethyl group, vinyl group, ethynyl group, methoxy group, difluoromethoxy group, trifluoromethoxy group, methylthio group, methylsulfinyl group and methylsulfonyl group,
- (4). a dibenzocycloheptene compound wherein R^1 of the compound represented by the formula (I) is selected from the group consisting of a hydrogen atom, a fluorine atom, a chlorine atom, a cyano group, trifluoromethyl group, hydroxymethyl group, 1-hydroxy-1-methylethyl group, ethynyl

group, methoxy group, difluoromethoxy group, trifluoromethoxy group, methylsulfinyl group and methylsulfonyl group,

5 (5). a dibenzocycloheptene compound wherein R^2 of the compound represented by the formula (I) is selected from the group consisting of a hydrogen atom, a fluorine atom, a chlorine atom, bromine atom, a nitro group, a cyano group, methyl group, ethyl group, propyl group, isopropyl group, methoxy group, ethoxy group, propoxy group and isopropoxy group,
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(6). a dibenzocycloheptene compound wherein R^2 of the compound represented by the formula (I) is selected from the group consisting of a hydrogen atom, a fluorine atom, a chlorine atom, a nitro group, a cyano group, methyl group, ethyl group, methoxy group and ethoxy group,
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(7). a dibenzocycloheptene compound wherein R^2 of the compound represented by the formula (I) is selected from the group consisting of a hydrogen atom, a fluorine atom, a chlorine atom, methyl group and methoxy group,

20 (8). a dibenzocycloheptene compound wherein R^2 of the compound represented by the formula (I) is a hydrogen atom,

(9). a dibenzocycloheptene compound wherein A of the compound represented by the formula (I) is selected from the group consisting of 2-thiazolyl, 1,3,4-thiadiazol-2-yl, 2-pyridyl, 2-pyrimidinyl, 2-benzoxazolyl, 2-benzothiazolyl, quinolin-2-yl, quinazolin-2-yl, 4-methyl-2-thiazolyl, 4-isopropyl-2-thiazolyl, 4-t-butyl-2-thiazolyl, 4-trifluoromethyl-2-thiazolyl, 5-methyl-1,3,4-thiadiazol-2-yl, 5-isopropyl-1,3,4-thiadiazol-2-yl, 5-t-butyl-1,3,4-thiadiazol-2-yl, 5-trifluoromethyl-1,3,4-thiadiazol-2-yl, 5,6-difluoro-2-pyridyl, 5,6-dichloro-2-pyridyl, 5,6-dimethyl-2-pyridyl, 5H-6,7-dihydrocyclopenta[b]pyridin-2-yl, 5,6,7,8-tetrahydroquinolin-2-yl, 5,6-difluoro-2-pyrimidinyl, 5,6-dichloro-2-pyrimidinyl, 5,6-dimethyl-2-pyrimidinyl, 6-trifluoromethyl-2-pyrimidinyl, 5H-6,7-dihydrocyclopenta[d]pyrimidin-2-yl, 5,6,7,8-tetrahydroquinazolin-2-yl, 6-
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fluoro-2-benzoxazolyl, 5-fluoro-2-benzoxazolyl, 5,6-
 difluoro-2-benzoxazolyl, 6-chloro-2-benzoxazolyl, 5-chloro-
 2-benzoxazolyl, 5,6-dichloro-2-benzoxazolyl, 5-chloro-6-
 fluoro-2-benzoxazolyl, 5-methyl-2-benzoxazolyl, 5-cyano-2-
 5 benzoxazolyl, 5-trifluoromethyl-2-benzoxazolyl, 5-methyl-
 thio-2-benzoxazolyl, 6-fluoro-2-benzothiazolyl, 5-fluoro-2-
 benzothiazolyl, 5,6-difluoro-2-benzothiazolyl, 6-chloro-2-
 benzothiazolyl, 5-chloro-2-benzothiazolyl, 5,6-dichloro-2-
 benzothiazolyl, 5-chloro-6-fluoro-2-benzothiazolyl, 5-
 10 methyl-2-benzothiazolyl, 5-cyano-2-benzothiazolyl, 5-tri-
 fluoromethyl-2-benzothiazolyl, 5-methylthio-2-benzothia-
 zolyl, 5-fluoroquinolin-2-yl, 6-fluoroquinolin-2-yl, 7-
 fluoroquinolin-2-yl, 5-chloroquinolin-2-yl, 6-chloro-
 quinolin-2-yl, 7-chloroquinolin-2-yl, 7-methylquinolin-2-
 15 yl, 7-trifluoromethylquinolin-2-yl, 7-methoxyquinolin-2-yl,
 7-difluoromethoxyquinolin-2-yl, 7-trifluoromethoxyquinolin-
 2-yl, 5,7-difluoroquinolin-2-yl, 6,7-difluoroquinolin-2-yl,
 5,7-dichloroquinolin-2-yl, 6,7-dichloroquinolin-2-yl, 5-
 chloro-7-fluoroquinolin-2-yl, 6-chloro-7-fluoroquinolin-2-
 20 yl, 7-chloro-5-fluoroquinolin-2-yl, 7-chloro-6-fluoro-
 quinolin-2-yl, 7-chloro-6-cyanoquinolin-2-yl, 7-cyano-6-
 fluoroquinolin-2-yl, 6-fluoro-7-trifluoromethylquinolin-2-
 yl, 5,6,7-trifluoroquinolin-2-yl, 5-fluoroquinazolin-2-yl,
 6-fluoroquinazolin-2-yl, 7-fluoroquinazolin-2-yl, 5-chloro-
 25 quinazolin-2-yl, 6-chloroquinazolin-2-yl, 7-chloroquina-
 zolin-2-yl, 7-methylquinazolin-2-yl, 7-trifluoromethyl-
 quinazolin-2-yl, 7-methoxyquinazolin-2-yl, 7-difluoro-
 methoxyquinazolin-2-yl, 7-trifluoromethoxyquinazolin-2-yl,
 5,7-difluoroquinazolin-2-yl, 6,7-difluoroquinazolin-2-yl,
 30 5,7-dichloroquinazolin-2-yl, 6,7-dichloroquinazolin-2-yl,
 5-chloro-7-fluoroquinazolin-2-yl, 6-chloro-7-fluoroquina-
 zolin-2-yl, 7-chloro-5-fluoroquinazolin-2-yl, 7-chloro-6-
 fluoroquinazolin-2-yl, 7-chloro-6-cyanoquinazolin-2-yl, 7-
 cyano-6-fluoroquinazolin-2-yl, 6-fluoro-7-trifluoromethyl-
 35 quinazolin-2-yl and 5,6,7-trifluoroquinazolin-2-yl group,
 (10). a dibenzocycloheptene compound wherein A of the

compound represented by the formula (I) is selected from the group consisting of 2-pyridyl, 2-benzothiazolyl, quinolin-2-yl, 5,6-difluoro-2-pyridyl, 5,6-dichloro-2-pyridyl, 5,6-dimethyl-2-pyridyl, 5,6,7,8-tetrahydro-quinolin-2-yl, 6-fluoro-2-benzothiazolyl, 5-fluoro-2-benzothiazolyl, 5,6-difluoro-2-benzothiazolyl, 6-chloro-2-benzothiazolyl, 5-chloro-2-benzothiazolyl, 5,6-dichloro-2-benzothiazolyl, 5-chloro-6-fluoro-2-benzothiazolyl, 5-methyl-2-benzothiazolyl, 5-cyano-2-benzothiazolyl, 5-trifluoromethyl-2-benzothiazolyl, 5-methylthio-2-benzothiazolyl, 5-fluoroquinolin-2-yl, 6-fluoroquinolin-2-yl, 7-fluoroquinolin-2-yl, 5-chloroquinolin-2-yl, 6-chloroquinolin-2-yl, 7-chloroquinolin-2-yl, 7-methylquinolin-2-yl, 7-trifluoromethylquinolin-2-yl, 7-methoxyquinolin-2-yl, 7-difluoromethoxyquinolin-2-yl, 7-trifluoromethoxyquinolin-2-yl, 5,7-difluoroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 5,7-dichloroquinolin-2-yl, 6,7-dichloroquinolin-2-yl, 5-chloro-7-fluoroquinolin-2-yl, 6-chloro-7-fluoroquinolin-2-yl, 7-chloro-5-fluoroquinolin-2-yl, 7-chloro-6-fluoroquinolin-2-yl, 7-chloro-6-cyanoquinolin-2-yl, 7-cyano-6-fluoroquinolin-2-yl, 6-fluoro-7-trifluoromethylquinolin-2-yl and 5,6,7-trifluoroquinolin-2-yl group,

(11). a dibenzocycloheptene compound wherein A of the compound represented by the formula (I) is selected from the group consisting of 5,6,7,8-tetrahydroquinolin-2-yl, 7-fluoroquinolin-2-yl, 7-chloroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 6,7-dichloroquinolin-2-yl and 7-chloro-6-fluoroquinolin-2-yl group,

(12). a dibenzocycloheptene compound wherein B of the compound represented by the formula (I) is a formula: -CH=CH-, a formula: -OCH₂- or a formula: -CH₂O-,

(13). a dibenzocycloheptene compound wherein X of the compound represented by the formula (I) is an oxygen atom or a sulfur atom,

(14). a dibenzocycloheptene compound wherein Y of the compound represented by the formula (I) is selected from

the group consisting of methylene, ethylene, trimethylene, tetramethylene, pentamethylene, fluoromethylene, difluoromethylene, 1-fluoroethylene, 2-fluoroethylene, 1,1-difluoroethylene, 2,2-difluoroethylene, ethylidene, 1-methylethylene, 2-methylethylene, 1,1-dimethylethylene, 2,2-dimethylethylene, 1-ethylethylene, 2-ethylethylene, 1-methoxyethylene, 2-methoxyethylene, 1-fluorotrimethylene, 2-fluorotrimethylene, 3-fluorotrimethylene, 1,1-difluorotrimethylene, 2,2-difluorotrimethylene, 3,3-difluorotrimethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethyltrimethylene, 2,2-dimethyltrimethylene, 3,3-dimethyltrimethylene, 2,2-diethyltrimethylene, 2-methoxytrimethylene, 3-methoxytrimethylene, 2,2-dimethoxytrimethylene, 3,3-dimethoxytrimethylene, (a-1) group, (a-2) group, (a-3) group, (a-4) group, (a-5) group and (a-6) group, (15). a dibenzocycloheptene compound wherein Y of the compound represented by the formula (I) is selected from the group consisting of methylene, ethylene, trimethylene, fluoromethylene, difluoromethylene, 1-fluoroethylene, 2-fluoroethylene, 1,1-difluoroethylene, 2,2-difluoroethylene, ethylidene, 1-methylethylene, 2-methylethylene, 1-fluorotrimethylene, 2-fluorotrimethylene, 3-fluorotrimethylene, 1,1-difluorotrimethylene, 2,2-difluorotrimethylene, 3,3-difluorotrimethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethyltrimethylene, 2,2-dimethyltrimethylene, 3,3-dimethyltrimethylene, (a-4) group, (a-5) group and (a-6) group, (16). a dibenzocycloheptene compound wherein Y of the compound represented by the formula (I) is selected from the group consisting of methylene, ethylene, trimethylene, difluoromethylene, 1-fluoroethylene, 2-fluoroethylene, 1,1-difluoroethylene, 2,2-difluoroethylene, ethylidene, 1-methylethylene, 2-methylethylene, 2,2-difluorotrimethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethyltrimethylene, 2,2-dimethyltrimethylene, 3,3-dimethyltrimethylene and (a-5) group,

(17). a dibenzocycloheptene compound wherein Y of the compound represented by the formula (I) is selected from the group consisting of methylene, ethylene, trimethylene, ethylidene, 1-methylethylene, 2-methylethylene and (a-5) group,

(18). a dibenzocycloheptene compound wherein Z of the compound represented by the formula (I) is selected from the group consisting of carboxy, 1H-tetrazol-5-yl, methane-sulfonylamino, ethanesulfonylamino, trifluoromethane-sulfonylamino, phenylsulfonylamino, p-fluorophenylsulfonylamino, p-chlorophenylsulfonylamino, o-methylphenylsulfonylamino, p-methylphenylsulfonylamino, p-trifluoromethylphenylsulfonylamino, o-methoxyphenylsulfonylamino, p-methoxyphenylsulfonylamino, p-difluoromethoxyphenylsulfonylamino, p-trifluoromethoxyphenylsulfonylamino, p-nitrophenylsulfonylamino, p-cyanophenylsulfonylamino, methane-sulfonylaminocarbonyl, ethanesulfonylaminocarbonyl, trifluoromethanesulfonylaminocarbonyl, phenylsulfonylaminocarbonyl, p-fluorophenylsulfonylaminocarbonyl, p-chlorophenylsulfonylaminocarbonyl, o-methylphenylsulfonylaminocarbonyl, p-methylphenylsulfonylaminocarbonyl, p-trifluoromethylphenylsulfonylaminocarbonyl, o-methoxyphenylsulfonylaminocarbonyl, p-methoxyphenylsulfonylaminocarbonyl, p-difluoromethoxyphenylsulfonylaminocarbonyl, p-trifluoromethoxyphenylsulfonylaminocarbonyl, p-nitrophenylsulfonylaminocarbonyl and p-cyanophenylsulfonylaminocarbonyl group,

(19). a dibenzocycloheptene compound wherein Z of the compound represented by the formula (I) is selected from the group consisting of carboxy, 1H-tetrazol-5-yl, methane-sulfonylamino, trifluoromethanesulfonylamino, phenylsulfonylamino, o-methylphenylsulfonylamino, p-methylphenylsulfonylamino, methanesulfonylaminocarbonyl, trifluoromethanesulfonylaminocarbonyl, phenylsulfonylaminocarbonyl, o-methylphenylsulfonylaminocarbonyl and p-methylphenylsulfonylaminocarbonyl group,

(20). a dibenzocycloheptene compound wherein Z of the

- compound represented by the formula (I) is selected from the group consisting of carboxy, methanesulfonylamino, trifluoromethanesulfonylamino, methanesulfonylaminocarbonyl and trifluoromethanesulfonylaminocarbonyl group,
- 5 (21). a dibenzocycloheptene compound wherein Z of the compound represented by the formula (I) is carboxyl group,
- (22). a dibenzocycloheptene compound wherein when Z of the compound represented by the formula (I) is carboxyl group, the protective group is selected from the group consisting
- 10 of C₁-C₄ alkyl group; benzyl group; C₁-C₂ alkyl group substituted by C₂-C₅ alkanoyloxy group; C₁-C₂ alkyl group substituted by (C₁-C₄ alkoxy)carbonyloxy group; and (5-methyl-2-oxo-1, 3-dioxolen-4-yl)methyl group,
- (23). a dibenzocycloheptene compound wherein when Z of the
- 15 compound represented by the formula (I) is carboxyl group, the protective group is selected from the group consisting of methyl, ethyl, propyl, isopropyl, butyl, isobutyl, acetoxymethyl, 1-acetoxyethyl, pivaloyloxymethyl and 1-pivaloyloxyethyl group,
- 20 (24). a dibenzocycloheptene compound wherein m of the compound represented by the formula (I) is 1, 2 or 3,
- (25). a dibenzocycloheptene compound wherein m of the compound represented by the formula (I) is 1 or 2,
- (26). a dibenzocycloheptene compound wherein n of the
- 25 compound represented by the formula (I) is 1 or 2,
- (27). a dibenzocycloheptene compound wherein n of the compound represented by the formula (I) is 1,
- and with respect to R¹, the preferable order is increased in the order of (1) to (4), with respect to R², the
- 30 preferable order is increased in the order of (5) to (8), with respect to A, the preferable order is increased in the order of (9) to (11), with respect to Y, the preferable order is increased in the order of (14) to (17), with respect to Z, the preferable order is increased in the
- 35 order of (18) to (21), with respect to the protective group when Z is a carboxyl group, the preferable order is

increased in the order of (22) to (23), with respect to m, the preferable order is increased in the order of (24) to (25), and with respect to n, the preferable order is increased in the order of (26) to (27).

5 Also, as the dibenzocycloheptene compound having the above-mentioned formula (I), a dibenzocycloheptene compound comprising a combination of two or more of the above-mentioned (1) to (27) is preferred.

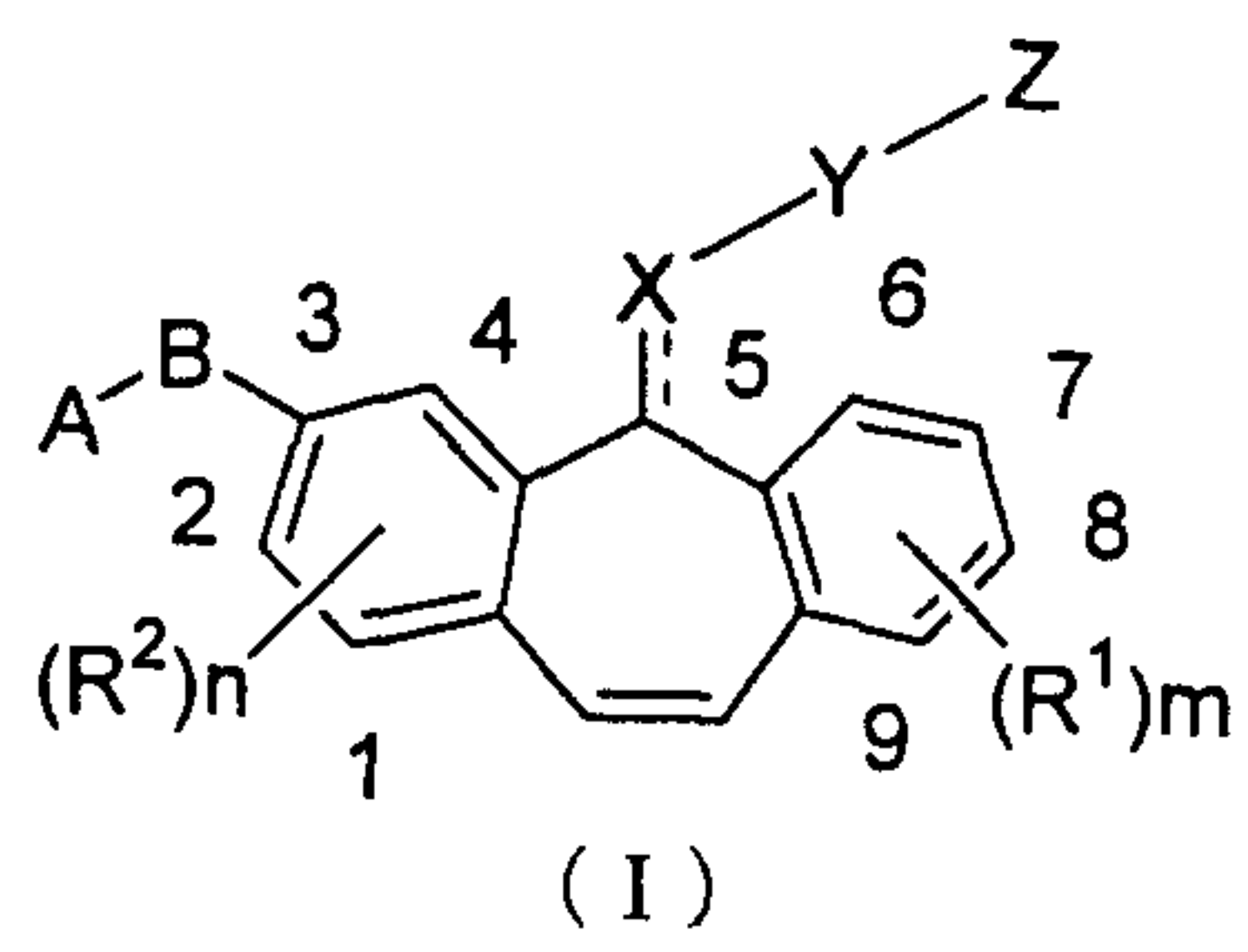
For example, the following may be mentioned,
 10 (28). a compound wherein R^1 of the compound represented by the formula (I) is selected from the group consisting of hydrogen atom, fluorine atom, chlorine atom, nitro group, cyano group, formyl group, 1H-tetrazol-5-yl group, methyl group, difluoromethyl group, trifluoromethyl group,
 15 hydroxymethyl group, 1-hydroxy-1-methylethyl group, vinyl group, ethynyl group, methoxy group, difluoromethoxy group, trifluoromethoxy group, methylthio group, methylsulfinyl group and methylsulfonyl group, R^2 is selected from the group consisting of hydrogen atom, fluorine atom, chlorine
 20 atom, methyl group and methoxy group, A is selected from the group consisting of 5,6,7,8-tetrahydroquinolin-2-yl, 7-fluoroquinolin-2-yl, 7-chloroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 6,7-dichloroquinolin-2-yl and 7-chloro-6-fluoroquinolin-2-yl group, B is a formula: $-CH=CH-$, a
 25 formula: $-OCH_2-$ or a formula: $-CH_2O-$, X is selected from the group consisting of an oxygen atom or a sulfur atom, Y is selected from the group consisting of methylene, ethylene, trimethylene, difluoromethylene, 1-fluoroethylene, 2-fluoroethylene, 1,1-difluoroethylene, 2,2-difluoroethylene,
 30 ethylidene, 1-methylethylene, 2-methylethylene, 2,2-difluorotrimethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethyltrimethylene, 2,2-dimethyltrimethylene, 3,3-dimethyltrimethylene and (a-5) group, Z is selected from the group consisting of carboxy, methanesulfonyl-
 35 amino, trifluoromethanesulfonylamino, methanesulfonylamino-carbonyl and trifluoromethanesulfonylaminocarbonyl group, m

is 1 or 2, and n is 1.

(29). a compound wherein R^1 of the compound represented by the formula (I) is selected from the group consisting of hydrogen atom, fluorine atom, chlorine atom, cyano group, trifluoromethyl group, hydroxymethyl group, 1-hydroxy-1-methylethyl group, ethynyl group, methoxy group, difluoromethoxy group, trifluoromethoxy group, methylsulfinyl group and methylsulfonyl group, R^2 is a hydrogen atom, A is selected from the group consisting of 5,6,7,8-tetrahydroquinolin-2-yl, 7-fluoroquinolin-2-yl, 7-chloroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 6,7-dichloroquinolin-2-yl and 7-chloro-6-fluoroquinolin-2-yl group, B is a formula: $-CH=CH-$, a formula: $-OCH_2-$ or a formula: $-CH_2O-$, X is an oxygen atom or a sulfur atom, Y is selected from the group consisting of methylene, ethylene, trimethylene, ethylidene, 1-methylethylene, 2-methylethylene and (a-5) group, Z is carboxy group, m is 1 or 2, and n is 1.

As preferred compounds in the compound (I), compounds in the following Table 1 can be specifically exemplified.

Table 1



No.	A	B	(R ²) _n	(R ¹) _m	X-Y-Z
1	6,7-diF-Q	-CH=CH-	H	H	-OCH ₂ COOH
2	6,7-diF-Q	-CH=CH-	H	H	-OCH(CH ₃)COOH
3	6,7-diF-Q	-CH=CH-	H	H	-OCH ₂ CH ₂ COOH
4	6,7-diF-Q	-CH=CH-	H	H	-OCH ₂ CH(CH ₃)COOH
5	6,7-diF-Q	-CH=CH-	H	H	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
6	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ COOH
7	6,7-diF-Q	-CH=CH-	H	H	-SCH(CH ₃)COOH
8	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
9	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)COOH
10	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₃) ₂ COOH
11	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₂ CH ₃)COOH
12	6,7-diF-Q	-CH=CH-	H	H	-SCH(CH ₃)CH ₂ COOH
13	6,7-diF-Q	-CH=CH-	H	H	-SC(CH ₃) ₂ CH ₂ COOH
14	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CH ₂ COOH
15	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)CH ₂ COOH
16	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₃) ₂ CH ₂ COOH
17	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
18	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CF ₂ COOH
19	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CF ₂ CH ₂ COOH
20	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ -Tet
21	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ -Tet
22	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ NHSO ₂ CF ₃
23	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ CH ₃
24	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ CF ₃
25	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ Ph
26	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ (2-CH ₃ -Ph)
27	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ NHSO ₂ CF ₃
28	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CH ₃
29	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CF ₃
30	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ Ph
31	6,7-diF-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ (2-CH ₃ -Ph)
32	6,7-diF-Q	-CH=CH-	H	H	-CH ₂ COOH
33	6,7-diF-Q	-CH=CH-	H	H	-CH ₂ CH ₂ COOH

34	6,7-diF-Q	-CH=CH-	H	H	-CH ₂ CH ₂ CH ₂ COOH
35	6,7-diF-Q	-CH=CH-	H	9-F	-OCH ₂ COOH
36	6,7-diF-Q	-CH=CH-	H	9-F	-OCH(CH ₃)COOH
37	6,7-diF-Q	-CH=CH-	H	9-F	-OCH ₂ CH ₂ COOH
38	6,7-diF-Q	-CH=CH-	H	9-F	-SCH ₂ COOH
39	6,7-diF-Q	-CH=CH-	H	9-F	-SCH(CH ₃)COOH
40	6,7-diF-Q	-CH=CH-	H	9-F	-SCH ₂ CH ₂ COOH
41	6,7-diF-Q	-CH=CH-	H	9-F	-SCH ₂ CH(CH ₃)COOH
42	6,7-diF-Q	-CH=CH-	H	9-F	-SCH ₂ CH(CH ₂ CH ₃)COOH
43	6,7-diF-Q	-CH=CH-	H	9-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
44	6,7-diF-Q	-CH=CH-	H	8-F	-OCH ₂ COOH
45	6,7-diF-Q	-CH=CH-	H	8-F	-OCH ₂ CH ₂ COOH
46	6,7-diF-Q	-CH=CH-	H	8-F	-SCH ₂ COOH
47	6,7-diF-Q	-CH=CH-	H	8-F	-SCH ₂ CH ₂ COOH
48	6,7-diF-Q	-CH=CH-	H	8-F	-SCH ₂ CH(CH ₃)COOH
49	6,7-diF-Q	-CH=CH-	H	8-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
50	6,7-diF-Q	-CH=CH-	H	7-F	-OCH ₂ COOH
51	6,7-diF-Q	-CH=CH-	H	7-F	-OCH ₂ CH ₂ COOH
52	6,7-diF-Q	-CH=CH-	H	7-F	-SCH ₂ COOH
53	6,7-diF-Q	-CH=CH-	H	7-F	-SCH ₂ CH ₂ COOH
54	6,7-diF-Q	-CH=CH-	H	7-F	-SCH ₂ CH(CH ₃)COOH
55	6,7-diF-Q	-CH=CH-	H	6-F	-OCH ₂ COOH
56	6,7-diF-Q	-CH=CH-	H	6-F	-OCH(CH ₃)COOH
57	6,7-diF-Q	-CH=CH-	H	6-F	-OCH ₂ CH ₂ COOH
58	6,7-diF-Q	-CH=CH-	H	6-F	-OCH ₂ CH(CH ₃)COOH
59	6,7-diF-Q	-CH=CH-	H	6-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
60	6,7-diF-Q	-CH=CH-	H	6-F	-SCH ₂ COOH
61	6,7-diF-Q	-CH=CH-	H	6-F	-SCH(CH ₃)COOH
62	6,7-diF-Q	-CH=CH-	H	6-F	-SCH ₂ CH ₂ COOH
63	6,7-diF-Q	-CH=CH-	H	6-F	-SCH ₂ CH(CH ₃)COOH
64	6,7-diF-Q	-CH=CH-	H	6-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
65	6,7-diF-Q	-CH=CH-	H	9-Cl	-OCH ₂ COOH
66	6,7-diF-Q	-CH=CH-	H	9-Cl	-OCH(CH ₃)COOH

67	6,7-diF-Q	-CH=CH-	H	9-Cl	-OCH ₂ CH ₂ COOH
68	6,7-diF-Q	-CH=CH-	H	9-Cl	-OCH ₂ CH(CH ₃)COOH
69	6,7-diF-Q	-CH=CH-	H	9-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
70	6,7-diF-Q	-CH=CH-	H	9-Cl	-SCH ₂ COOH
71	6,7-diF-Q	-CH=CH-	H	9-Cl	-SCH(CH ₃)COOH
72	6,7-diF-Q	-CH=CH-	H	9-Cl	-SCH ₂ CH ₂ COOH
73	6,7-diF-Q	-CH=CH-	H	9-Cl	-SCH ₂ CH(CH ₃)COOH
74	6,7-diF-Q	-CH=CH-	H	9-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
75	6,7-diF-Q	-CH=CH-	H	8-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
76	6,7-diF-Q	-CH=CH-	H	8-Cl	-OCH ₂ COOH
77	6,7-diF-Q	-CH=CH-	H	8-Cl	-OCH ₂ CH ₂ COOH
78	6,7-diF-Q	-CH=CH-	H	8-Cl	-SCH ₂ COOH
79	6,7-diF-Q	-CH=CH-	H	8-Cl	-SCH ₂ CH ₂ COOH
80	6,7-diF-Q	-CH=CH-	H	8-Cl	-SCH ₂ CH(CH ₃)COOH
81	6,7-diF-Q	-CH=CH-	H	7-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
82	6,7-diF-Q	-CH=CH-	H	7-Cl	-OCH ₂ COOH
83	6,7-diF-Q	-CH=CH-	H	7-Cl	-OCH ₂ CH ₂ COOH
84	6,7-diF-Q	-CH=CH-	H	7-Cl	-SCH ₂ COOH
85	6,7-diF-Q	-CH=CH-	H	7-Cl	-SCH ₂ CH ₂ COOH
86	6,7-diF-Q	-CH=CH-	H	7-Cl	-SCH ₂ CH(CH ₃)COOH
87	6,7-diF-Q	-CH=CH-	H	6-Cl	-OCH ₂ COOH
88	6,7-diF-Q	-CH=CH-	H	6-Cl	-OCH(CH ₃)COOH
89	6,7-diF-Q	-CH=CH-	H	6-Cl	-OCH ₂ CH ₂ COOH
90	6,7-diF-Q	-CH=CH-	H	6-Cl	-OCH ₂ CH(CH ₃)COOH
91	6,7-diF-Q	-CH=CH-	H	6-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
92	6,7-diF-Q	-CH=CH-	H	6-Cl	-SCH ₂ COOH
93	6,7-diF-Q	-CH=CH-	H	6-Cl	-SCH(CH ₃)COOH
94	6,7-diF-Q	-CH=CH-	H	6-Cl	-SCH ₂ CH ₂ COOH
95	6,7-diF-Q	-CH=CH-	H	6-Cl	-SCH ₂ CH(CH ₃)COOH
96	6,7-diF-Q	-CH=CH-	H	6-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
97	6,7-diF-Q	-CH=CH-	H	9-CN	-OCH ₂ COOH
98	6,7-diF-Q	-CH=CH-	H	9-CN	-OCH(CH ₃)COOH
99	6,7-diF-Q	-CH=CH-	H	9-CN	-OCH ₂ CH ₂ COOH

100	6,7-diF-Q	-CH=CH-	H	9-CN	-OCH ₂ CH(CH ₃)COOH
101	6,7-diF-Q	-CH=CH-	H	9-CN	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
102	6,7-diF-Q	-CH=CH-	H	9-CN	-SCH ₂ COOH
103	6,7-diF-Q	-CH=CH-	H	9-CN	-SCH(CH ₃)COOH
104	6,7-diF-Q	-CH=CH-	H	9-CN	-SCH ₂ CH ₂ COOH
105	6,7-diF-Q	-CH=CH-	H	9-CN	-SCH ₂ CH(CH ₃)COOH
106	6,7-diF-Q	-CH=CH-	H	9-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
107	6,7-diF-Q	-CH=CH-	H	8-CN	-OCH ₂ COOH
108	6,7-diF-Q	-CH=CH-	H	8-CN	-OCH ₂ CH ₂ COOH
109	6,7-diF-Q	-CH=CH-	H	8-CN	-SCH ₂ COOH
110	6,7-diF-Q	-CH=CH-	H	8-CN	-SCH ₂ CH ₂ COOH
111	6,7-diF-Q	-CH=CH-	H	8-CN	-SCH ₂ CH(CH ₃)COOH
112	6,7-diF-Q	-CH=CH-	H	8-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
113	6,7-diF-Q	-CH=CH-	H	7-CN	-OCH ₂ COOH
114	6,7-diF-Q	-CH=CH-	H	7-CN	-OCH ₂ CH ₂ COOH
115	6,7-diF-Q	-CH=CH-	H	7-CN	-SCH ₂ COOH
116	6,7-diF-Q	-CH=CH-	H	7-CN	-SCH ₂ CH ₂ COOH
117	6,7-diF-Q	-CH=CH-	H	7-CN	-SCH ₂ CH(CH ₃)COOH
118	6,7-diF-Q	-CH=CH-	H	9-CH ₃	-OCH ₂ COOH
119	6,7-diF-Q	-CH=CH-	H	9-CH ₃	-OCH ₂ CH ₂ COOH
120	6,7-diF-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ COOH
121	6,7-diF-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH ₂ COOH
122	6,7-diF-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH(CH ₃)COOH
123	6,7-diF-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH(CH ₂ CH ₃)COOH
124	6,7-diF-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
125	6,7-diF-Q	-CH=CH-	H	8-CH ₃	-OCH ₂ COOH
126	6,7-diF-Q	-CH=CH-	H	8-CH ₃	-OCH ₂ CH ₂ COOH
127	6,7-diF-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ COOH
128	6,7-diF-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ CH ₂ COOH
129	6,7-diF-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ CH(CH ₃)COOH
130	6,7-diF-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
131	6,7-diF-Q	-CH=CH-	-H	7-CH ₃	-OCH ₂ COOH
132	6,7-diF-Q	-CH=CH-	H	7-CH ₃	-OCH ₂ CH ₂ COOH

133	6,7-diF-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ COOH
134	6,7-diF-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ CH ₂ COOH
135	6,7-diF-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ CH(CH ₃)COOH
136	6,7-diF-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ COOH
137	6,7-diF-Q	-CH=CH-	H	9-CF ₃	-OCH(CH ₃)COOH
138	6,7-diF-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ CH ₂ COOH
139	6,7-diF-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ CH(CH ₃)COOH
140	6,7-diF-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
141	6,7-diF-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ COOH
142	6,7-diF-Q	-CH=CH-	H	9-CF ₃	-SCH(CH ₃)COOH
143	6,7-diF-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ CH ₂ COOH
144	6,7-diF-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ CH(CH ₃)COOH
145	6,7-diF-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
146	6,7-diF-Q	-CH=CH-	H	8-CF ₃	-OCH ₂ COOH
147	6,7-diF-Q	-CH=CH-	H	8-CF ₃	-OCH ₂ CH ₂ COOH
148	6,7-diF-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ COOH
149	6,7-diF-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ CH ₂ COOH
150	6,7-diF-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ CH(CH ₃)COOH
151	6,7-diF-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
152	6,7-diF-Q	-CH=CH-	H	7-CF ₃	-OCH ₂ COOH
153	6,7-diF-Q	-CH=CH-	H	7-CF ₃	-OCH ₂ CH ₂ COOH
154	6,7-diF-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ COOH
155	6,7-diF-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ CH ₂ COOH
156	6,7-diF-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ CH(CH ₃)COOH
157	6,7-diF-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ COOH
158	6,7-diF-Q	-CH=CH-	H	9-C≡CH	-OCH(CH ₃)COOH
159	6,7-diF-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ CH ₂ COOH
160	6,7-diF-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ CH(CH ₃)COOH
161	6,7-diF-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
162	6,7-diF-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ COOH
163	6,7-diF-Q	-CH=CH-	H	9-C≡CH	-SCH(CH ₃)COOH
164	6,7-diF-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ CH ₂ COOH
165	6,7-diF-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ CH(CH ₃)COOH

166	6,7-diF-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
167	6,7-diF-Q	-CH=CH-	H	8-C≡CH	-OCH ₂ COOH
168	6,7-diF-Q	-CH=CH-	H	8-C≡CH	-OCH ₂ CH ₂ COOH
169	6,7-diF-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ COOH
170	6,7-diF-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ CH ₂ COOH
171	6,7-diF-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ CH(CH ₃)COOH
172	6,7-diF-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
173	6,7-diF-Q	-CH=CH-	H	7-C≡CH	-OCH ₂ COOH
174	6,7-diF-Q	-CH=CH-	H	7-C≡CH	-OCH ₂ CH ₂ COOH
175	6,7-diF-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ COOH
176	6,7-diF-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ CH ₂ COOH
177	6,7-diF-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ CH(CH ₃)COOH
178	6,7-diF-Q	-CH=CH-	H	9-CH ₂ OH	-OCH ₂ COOH
179	6,7-diF-Q	-CH=CH-	H	9-CH ₂ OH	-SCH ₂ CH ₂ COOH
180	6,7-diF-Q	-CH=CH-	H	8-CH ₂ OH	-OCH ₂ COOH
181	6,7-diF-Q	-CH=CH-	H	8-CH ₂ OH	-SCH ₂ CH ₂ COOH
182	6,7-diF-Q	-CH=CH-	H	7-CH ₂ OH	-OCH ₂ COOH
183	6,7-diF-Q	-CH=CH-	H	7-CH ₂ OH	-SCH ₂ CH ₂ COOH
184	6,7-diF-Q	-CH=CH-	H	9- C(CH ₃) ₂ OH	-OCH ₂ COOH
185	6,7-diF-Q	-CH=CH-	H	9- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
186	6,7-diF-Q	-CH=CH-	H	8- C(CH ₃) ₂ OH	-OCH ₂ COOH
187	6,7-diF-Q	-CH=CH-	H	8- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
188	6,7-diF-Q	-CH=CH-	H	7- C(CH ₃) ₂ OH	-OCH ₂ COOH
189	6,7-diF-Q	-CH=CH-	H	7- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
190	6,7-diF-Q	-CH=CH-	H	9-OCH ₃	-OCH ₂ COOH
191	6,7-diF-Q	-CH=CH-	H	9-OCH ₃	-SCH ₂ CH ₂ COOH
192	6,7-diF-Q	-CH=CH-	H	8-OCH ₃	-OCH ₂ COOH

193	6,7-diF-Q	-CH=CH-	H	8-OCH ₃	-SCH ₂ CH ₂ COOH
194	6,7-diF-Q	-CH=CH-	H	7-OCH ₃	-OCH ₂ COOH
195	6,7-diF-Q	-CH=CH-	H	7-OCH ₃	-SCH ₂ CH ₂ COOH
196	6,7-diF-Q	-CH=CH-	H	9-OCHF ₂	-OCH ₂ COOH
197	6,7-diF-Q	-CH=CH-	H	9-OCHF ₂	-SCH ₂ CH ₂ COOH
198	6,7-diF-Q	-CH=CH-	H	8-OCHF ₂	-OCH ₂ COOH
199	6,7-diF-Q	-CH=CH-	H	8-OCHF ₂	-SCH ₂ CH ₂ COOH
200	6,7-diF-Q	-CH=CH-	H	7-OCHF ₂	-OCH ₂ COOH
201	6,7-diF-Q	-CH=CH-	H	7-OCHF ₂	-SCH ₂ CH ₂ COOH
202	6,7-diF-Q	-CH=CH-	H	9-OCF ₃	-OCH ₂ COOH
203	6,7-diF-Q	-CH=CH-	H	9-OCF ₃	-SCH ₂ CH ₂ COOH
204	6,7-diF-Q	-CH=CH-	H	8-OCF ₃	-OCH ₂ COOH
205	6,7-diF-Q	-CH=CH-	H	8-OCF ₃	-SCH ₂ CH ₂ COOH
206	6,7-diF-Q	-CH=CH-	H	7-OCF ₃	-OCH ₂ COOH
207	6,7-diF-Q	-CH=CH-	H	7-OCF ₃	-SCH ₂ CH ₂ COOH
208	6,7-diF-Q	-CH=CH-	H	9-SOCH ₃	-OCH ₂ COOH
209	6,7-diF-Q	-CH=CH-	H	9-SOCH ₃	-SCH ₂ CH ₂ COOH
210	6,7-diF-Q	-CH=CH-	H	8-SOCH ₃	-OCH ₂ COOH
211	6,7-diF-Q	-CH=CH-	H	8-SOCH ₃	-SCH ₂ CH ₂ COOH
212	6,7-diF-Q	-CH=CH-	H	7-SOCH ₃	-OCH ₂ COOH
213	6,7-diF-Q	-CH=CH-	H	7-SOCH ₃	-SCH ₂ CH ₂ COOH
214	6,7-diF-Q	-CH=CH-	H	9-SO ₂ CH ₃	-OCH ₂ COOH
215	6,7-diF-Q	-CH=CH-	H	9-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
216	6,7-diF-Q	-CH=CH-	H	8-SO ₂ CH ₃	-OCH ₂ COOH
217	6,7-diF-Q	-CH=CH-	H	8-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
218	6,7-diF-Q	-CH=CH-	H	7-SO ₂ CH ₃	-OCH ₂ COOH
219	6,7-diF-Q	-CH=CH-	H	7-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
220	6,7-diF-Q	-CH=CH-	H	9-CH=CH ₂	-SCH ₂ CH ₂ COOH
221	6,7-diF-Q	-CH=CH-	H	8-CH=CH ₂	-SCH ₂ CH ₂ COOH
222	6,7-diF-Q	-CH=CH-	H	7-CH=CH ₂	-SCH ₂ CH ₂ COOH
223	6,7-diF-Q	-CH=CH-	H	9-NO ₂	-OCH ₂ COOH
224	6,7-diF-Q	-CH=CH-	H	9-NO ₂	-SCH ₂ CH ₂ COOH
225	6,7-diF-Q	-CH=CH-	H	8-NO ₂	-OCH ₂ COOH

226	6,7-diF-Q	-CH=CH-	H	8-NO ₂	-SCH ₂ CH ₂ COOH
227	6,7-diF-Q	-CH=CH-	H	7-NO ₂	-OCH ₂ COOH
228	6,7-diF-Q	-CH=CH-	H	7-NO ₂	-SCH ₂ CH ₂ COOH
229	6,7-diF-Q	-CH=CH-	1-F	H	-SCH ₂ CH ₂ COOH
230	6,7-diF-Q	-CH=CH-	2-F	H	-SCH ₂ CH ₂ COOH
231	6,7-diF-Q	-CH=CH-	1-Cl	H	-SCH ₂ CH ₂ COOH
232	6,7-diF-Q	-CH=CH-	2-Cl	H	-SCH ₂ CH ₂ COOH
233	6,7-diF-Q	-CH=CH-	1-CH ₃	H	-SCH ₂ CH ₂ COOH
234	6,7-diF-Q	-CH=CH-	2-CH ₃	H	-SCH ₂ CH ₂ COOH
235	6,7-diF-Q	-CH=CH-	1-OCH ₃	H	-SCH ₂ CH ₂ COOH
236	6,7-diF-Q	-CH=CH-	2-OCH ₃	H	-SCH ₂ CH ₂ COOH
237	6,7-diF-Q	-CH=CH-	1-NO ₂	H	-SCH ₂ CH ₂ COOH
238	6,7-diF-Q	-CH=CH-	2-NO ₂	H	-SCH ₂ CH ₂ COOH
239	6,7-diF-Q	-CH=CH-	1-CN	H	-SCH ₂ CH ₂ COOH
240	6,7-diF-Q	-CH=CH-	2-CN	H	-SCH ₂ CH ₂ COOH
241	6,7-diF-Q	-CH ₂ O-	H	H	-OCH ₂ COOH
242	6,7-diF-Q	-CH ₂ O-	H	H	-OCH(CH ₃)COOH
243	6,7-diF-Q	-CH ₂ O-	H	H	-OCH ₂ CH ₂ COOH
244	6,7-diF-Q	-CH ₂ O-	H	H	-OCH ₂ CH(CH ₃)COOH
245	6,7-diF-Q	-CH ₂ O-	H	H	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
246	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ COOH
247	6,7-diF-Q	-CH ₂ O-	H	H	-SCH(CH ₃)COOH
248	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
249	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)COOH
250	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₃) ₂ COOH
251	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₂ CH ₃)COOH
252	6,7-diF-Q	-CH ₂ O-	H	H	-SCH(CH ₃)CH ₂ COOH
253	6,7-diF-Q	-CH ₂ O-	H	H	-SC(CH ₃) ₂ CH ₂ COOH
254	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CH ₂ COOH
255	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)CH ₂ COOH
256	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₃) ₂ CH ₂ COOH
257	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
258	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CF ₂ COOH

259	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CF ₂ CH ₂ COOH
260	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ -Tet
261	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ -Tet
262	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ NHSO ₂ CF ₃
263	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ CH ₃
264	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ CF ₃
265	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ Ph
266	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ (2-CH ₃ -Ph)
267	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ NHSO ₂ CF ₃
268	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CH ₃
269	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CF ₃
270	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ Ph
271	6,7-diF-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ (2-CH ₃ -Ph)
272	6,7-diF-Q	-CH ₂ O-	H	H	-CH ₂ COOH
273	6,7-diF-Q	-CH ₂ O-	H	H	-CH ₂ CH ₂ COOH
274	6,7-diF-Q	-CH ₂ O-	H	H	-CH ₂ CH ₂ CH ₂ COOH
275	6,7-diF-Q	-CH ₂ O-	H	9-F	-OCH ₂ COOH
276	6,7-diF-Q	-CH ₂ O-	H	9-F	-OCH(CH ₃)COOH
277	6,7-diF-Q	-CH ₂ O-	H	9-F	-OCH ₂ CH ₂ COOH
278	6,7-diF-Q	-CH ₂ O-	H	9-F	-OCH ₂ CH(CH ₃)COOH
279	6,7-diF-Q	-CH ₂ O-	H	9-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
280	6,7-diF-Q	-CH ₂ O-	H	9-F	-SCH ₂ COOH
281	6,7-diF-Q	-CH ₂ O-	H	9-F	-SCH(CH ₃)COOH
282	6,7-diF-Q	-CH ₂ O-	H	9-F	-SCH ₂ CH ₂ COOH
283	6,7-diF-Q	-CH ₂ O-	H	9-F	-SCH ₂ CH(CH ₃)COOH
284	6,7-diF-Q	-CH ₂ O-	H	9-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
285	6,7-diF-Q	-CH ₂ O-	H	8-F	-OCH ₂ COOH
286	6,7-diF-Q	-CH ₂ O-	H	8-F	-OCH ₂ CH ₂ COOH
287	6,7-diF-Q	-CH ₂ O-	H	8-F	-SCH ₂ COOH
288	6,7-diF-Q	-CH ₂ O-	H	8-F	-SCH ₂ CH ₂ COOH
289	6,7-diF-Q	-CH ₂ O-	H	8-F	-SCH ₂ CH(CH ₃)COOH
290	6,7-diF-Q	-CH ₂ O-	H	8-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
291	6,7-diF-Q	-CH ₂ O-	H	7-F	-OCH ₂ COOH

292	6,7-diF-Q	-CH ₂ O-	H	7-F	-OCH ₂ CH ₂ COOH
293	6,7-diF-Q	-CH ₂ O-	H	7-F	-SCH ₂ COOH
294	6,7-diF-Q	-CH ₂ O-	H	7-F	-SCH ₂ CH ₂ COOH
295	6,7-diF-Q	-CH ₂ O-	H	7-F	-SCH ₂ CH(CH ₃)COOH
296	6,7-diF-Q	-CH ₂ O-	H	6-F	-OCH ₂ COOH
297	6,7-diF-Q	-CH ₂ O-	H	6-F	-OCH(CH ₃)COOH
298	6,7-diF-Q	-CH ₂ O-	H	6-F	-OCH ₂ CH ₂ COOH
299	6,7-diF-Q	-CH ₂ O-	H	6-F	-OCH ₂ CH(CH ₃)COOH
300	6,7-diF-Q	-CH ₂ O-	H	6-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
301	6,7-diF-Q	-CH ₂ O-	H	6-F	-SCH ₂ COOH
302	6,7-diF-Q	-CH ₂ O-	H	6-F	-SCH(CH ₃)COOH
303	6,7-diF-Q	-CH ₂ O-	H	6-F	-SCH ₂ CH ₂ COOH
304	6,7-diF-Q	-CH ₂ O-	H	6-F	-SCH ₂ CH(CH ₃)COOH
305	6,7-diF-Q	-CH ₂ O-	H	6-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
306	6,7-diF-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ COOH
307	6,7-diF-Q	-CH ₂ O-	H	9-Cl	-OCH(CH ₃)COOH
308	6,7-diF-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ CH ₂ COOH
309	6,7-diF-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ CH(CH ₃)COOH
310	6,7-diF-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
311	6,7-diF-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ COOH
312	6,7-diF-Q	-CH ₂ O-	H	9-Cl	-SCH(CH ₃)COOH
313	6,7-diF-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ CH ₂ COOH
314	6,7-diF-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ CH(CH ₃)COOH
315	6,7-diF-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
316	6,7-diF-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
317	6,7-diF-Q	-CH ₂ O-	H	8-Cl	-OCH ₂ COOH
318	6,7-diF-Q	-CH ₂ O-	H	8-Cl	-OCH ₂ CH ₂ COOH
319	6,7-diF-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ COOH
320	6,7-diF-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ CH ₂ COOH
321	6,7-diF-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ CH(CH ₃)COOH
322	6,7-diF-Q	-CH ₂ O-	H	7-Cl	-OCH ₂ COOH
323	6,7-diF-Q	-CH ₂ O-	H	7-Cl	-OCH ₂ CH ₂ COOH
324	6,7-diF-Q	-CH ₂ O-	H	7-Cl	-SCH ₂ COOH

325	6,7-diF-Q	-CH ₂ O-	H	7-Cl	-SCH ₂ CH ₂ COOH
326	6,7-diF-Q	-CH ₂ O-	H	7-Cl	-SCH ₂ CH(CH ₃)COOH
327	6,7-diF-Q	-CH ₂ O-	H	9-CN	-OCH ₂ COOH
328	6,7-diF-Q	-CH ₂ O-	H	9-CN	-OCH(CH ₃)COOH
329	6,7-diF-Q	-CH ₂ O-	H	9-CN	-OCH ₂ CH ₂ COOH
330	6,7-diF-Q	-CH ₂ O-	H	9-CN	-OCH ₂ CH(CH ₃)COOH
331	6,7-diF-Q	-CH ₂ O-	H	9-CN	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
332	6,7-diF-Q	-CH ₂ O-	H	9-CN	-SCH ₂ COOH
333	6,7-diF-Q	-CH ₂ O-	H	9-CN	-SCH(CH ₃)COOH
334	6,7-diF-Q	-CH ₂ O-	H	9-CN	-SCH ₂ CH ₂ COOH
335	6,7-diF-Q	-CH ₂ O-	H	9-CN	-SCH ₂ CH(CH ₃)COOH
336	6,7-diF-Q	-CH ₂ O-	H	9-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
337	6,7-diF-Q	-CH ₂ O-	H	8-CN	-OCH ₂ COOH
338	6,7-diF-Q	-CH ₂ O-	H	8-CN	-OCH ₂ CH ₂ COOH
339	6,7-diF-Q	-CH ₂ O-	H	8-CN	-SCH ₂ COOH
340	6,7-diF-Q	-CH ₂ O-	H	8-CN	-SCH ₂ CH ₂ COOH
341	6,7-diF-Q	-CH ₂ O-	H	8-CN	-SCH ₂ CH(CH ₃)COOH
342	6,7-diF-Q	-CH ₂ O-	H	8-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
343	6,7-diF-Q	-CH ₂ O-	H	7-CN	-OCH ₂ COOH
344	6,7-diF-Q	-CH ₂ O-	H	7-CN	-OCH ₂ CH ₂ COOH
345	6,7-diF-Q	-CH ₂ O-	H	7-CN	-SCH ₂ COOH
346	6,7-diF-Q	-CH ₂ O-	H	7-CN	-SCH ₂ CH ₂ COOH
347	6,7-diF-Q	-CH ₂ O-	H	7-CN	-SCH ₂ CH(CH ₃)COOH
348	6,7-diF-Q	-CH ₂ O-	H	9-CH ₃	-OCH ₂ COOH
349	6,7-diF-Q	-CH ₂ O-	H	9-CH ₃	-OCH ₂ CH ₂ COOH
350	6,7-diF-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ COOH
351	6,7-diF-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ CH ₂ COOH
352	6,7-diF-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ CH(CH ₃)COOH
353	6,7-diF-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ CH(CH ₂ CH ₃)COOH
354	6,7-diF-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
355	6,7-diF-Q	-CH ₂ O-	H	8-CH ₃	-OCH ₂ COOH
356	6,7-diF-Q	-CH ₂ O-	H	8-CH ₃	-OCH ₂ CH ₂ COOH
357	6,7-diF-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ COOH

358	6,7-diF-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ CH ₂ COOH
359	6,7-diF-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ CH(CH ₃)COOH
360	6,7-diF-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
361	6,7-diF-Q	-CH ₂ O-	H	7-CH ₃	-OCH ₂ COOH
362	6,7-diF-Q	-CH ₂ O-	H	7-CH ₃	-OCH ₂ CH ₂ COOH
363	6,7-diF-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ COOH
364	6,7-diF-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ CH ₂ COOH
365	6,7-diF-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ CH(CH ₃)COOH
366	6,7-diF-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ COOH
367	6,7-diF-Q	-CH ₂ O-	H	9-CF ₃	-OCH(CH ₃)COOH
368	6,7-diF-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ CH ₂ COOH
369	6,7-diF-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ CH(CH ₃)COOH
370	6,7-diF-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
371	6,7-diF-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ COOH
372	6,7-diF-Q	-CH ₂ O-	H	9-CF ₃	-SCH(CH ₃)COOH
373	6,7-diF-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ CH ₂ COOH
374	6,7-diF-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ CH(CH ₃)COOH
375	6,7-diF-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
376	6,7-diF-Q	-CH ₂ O-	H	8-CF ₃	-OCH ₂ COOH
377	6,7-diF-Q	-CH ₂ O-	H	8-CF ₃	-OCH ₂ CH ₂ COOH
378	6,7-diF-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ COOH
379	6,7-diF-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ CH ₂ COOH
380	6,7-diF-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ CH(CH ₃)COOH
381	6,7-diF-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
382	6,7-diF-Q	-CH ₂ O-	H	7-CF ₃	-OCH ₂ COOH
383	6,7-diF-Q	-CH ₂ O-	H	7-CF ₃	-OCH ₂ CH ₂ COOH
384	6,7-diF-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ COOH
385	6,7-diF-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ CH ₂ COOH
386	6,7-diF-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ CH(CH ₃)COOH
387	6,7-diF-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ COOH
388	6,7-diF-Q	-CH ₂ O-	H	9-C≡CH	-OCH(CH ₃)COOH
389	6,7-diF-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ CH ₂ COOH
390	6,7-diF-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ CH(CH ₃)COOH

391	6,7-diF-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
392	6,7-diF-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ COOH
393	6,7-diF-Q	-CH ₂ O-	H	9-C≡CH	-SCH(CH ₃)COOH
394	6,7-diF-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ CH ₂ COOH
395	6,7-diF-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ CH(CH ₃)COOH
396	6,7-diF-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
397	6,7-diF-Q	-CH ₂ O-	H	8-C≡CH	-OCH ₂ COOH
398	6,7-diF-Q	-CH ₂ O-	H	8-C≡CH	-OCH ₂ CH ₂ COOH
399	6,7-diF-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ COOH
400	6,7-diF-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ CH ₂ COOH
401	6,7-diF-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ CH(CH ₃)COOH
402	6,7-diF-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
403	6,7-diF-Q	-CH ₂ O-	H	7-C≡CH	-OCH ₂ COOH
404	6,7-diF-Q	-CH ₂ O-	H	7-C≡CH	-OCH ₂ CH ₂ COOH
405	6,7-diF-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ COOH
406	6,7-diF-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ CH ₂ COOH
407	6,7-diF-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ CH(CH ₃)COOH
408	6,7-diF-Q	-CH ₂ O-	H	9-CH ₂ OH	-OCH ₂ COOH
409	6,7-diF-Q	-CH ₂ O-	H	9-CH ₂ OH	-SCH ₂ CH ₂ COOH
410	6,7-diF-Q	-CH ₂ O-	H	8-CH ₂ OH	-OCH ₂ COOH
411	6,7-diF-Q	-CH ₂ O-	H	8-CH ₂ OH	-SCH ₂ CH ₂ COOH
412	6,7-diF-Q	-CH ₂ O-	H	7-CH ₂ OH	-OCH ₂ COOH
413	6,7-diF-Q	-CH ₂ O-	H	7-CH ₂ OH	-SCH ₂ CH ₂ COOH
414	6,7-diF-Q	-CH ₂ O-	H	9- C(CH ₃) ₂ OH	-OCH ₂ COOH
415	6,7-diF-Q	-CH ₂ O-	H	9- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
416	6,7-diF-Q	-CH ₂ O-	H	8- C(CH ₃) ₂ OH	-OCH ₂ COOH
417	6,7-diF-Q	-CH ₂ O-	H	8- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
418	6,7-diF-Q	-CH ₂ O-	H	7- C(CH ₃) ₂ OH	-OCH ₂ COOH

419	6,7-diF-Q	-CH ₂ O-	H	7- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
420	6,7-diF-Q	-CH ₂ O-	H	9-OCH ₃	-OCH ₂ COOH
421	6,7-diF-Q	-CH ₂ O-	H	9-OCH ₃	-SCH ₂ CH ₂ COOH
422	6,7-diF-Q	-CH ₂ O-	H	8-OCH ₃	-OCH ₂ COOH
423	6,7-diF-Q	-CH ₂ O-	H	8-OCH ₃	-SCH ₂ CH ₂ COOH
424	6,7-diF-Q	-CH ₂ O-	H	7-OCH ₃	-OCH ₂ COOH
425	6,7-diF-Q	-CH ₂ O-	H	7-OCH ₃	-SCH ₂ CH ₂ COOH
426	6,7-diF-Q	-CH ₂ O-	H	9-OCHF ₂	-OCH ₂ COOH
427	6,7-diF-Q	-CH ₂ O-	H	9-OCHF ₂	-SCH ₂ CH ₂ COOH
428	6,7-diF-Q	-CH ₂ O-	H	8-OCHF ₂	-OCH ₂ COOH
429	6,7-diF-Q	-CH ₂ O-	H	8-OCHF ₂	-SCH ₂ CH ₂ COOH
430	6,7-diF-Q	-CH ₂ O-	H	7-OCHF ₂	-OCH ₂ COOH
431	6,7-diF-Q	-CH ₂ O-	H	7-OCHF ₂	-SCH ₂ CH ₂ COOH
432	6,7-diF-Q	-CH ₂ O-	H	9-OCF ₃	-OCH ₂ COOH
433	6,7-diF-Q	-CH ₂ O-	H	9-OCF ₃	-SCH ₂ CH ₂ COOH
434	6,7-diF-Q	-CH ₂ O-	H	8-OCF ₃	-OCH ₂ COOH
435	6,7-diF-Q	-CH ₂ O-	H	8-OCF ₃	-SCH ₂ CH ₂ COOH
436	6,7-diF-Q	-CH ₂ O-	H	7-OCF ₃	-OCH ₂ COOH
437	6,7-diF-Q	-CH ₂ O-	H	7-OCF ₃	-SCH ₂ CH ₂ COOH
438	6,7-diF-Q	-CH ₂ O-	H	9-SOCH ₃	-OCH ₂ COOH
439	6,7-diF-Q	-CH ₂ O-	H	9-SOCH ₃	-SCH ₂ CH ₂ COOH
440	6,7-diF-Q	-CH ₂ O-	H	8-SOCH ₃	-OCH ₂ COOH
441	6,7-diF-Q	-CH ₂ O-	H	8-SOCH ₃	-SCH ₂ CH ₂ COOH
442	6,7-diF-Q	-CH ₂ O-	H	7-SOCH ₃	-OCH ₂ COOH
443	6,7-diF-Q	-CH ₂ O-	H	7-SOCH ₃	-SCH ₂ CH ₂ COOH
444	6,7-diF-Q	-CH ₂ O-	H	9-SO ₂ CH ₃	-OCH ₂ COOH
445	6,7-diF-Q	-CH ₂ O-	H	9-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
446	6,7-diF-Q	-CH ₂ O-	H	8-SO ₂ CH ₃	-OCH ₂ COOH
447	6,7-diF-Q	-CH ₂ O-	H	8-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
448	6,7-diF-Q	-CH ₂ O-	H	7-SO ₂ CH ₃	-OCH ₂ COOH
449	6,7-diF-Q	-CH ₂ O-	H	7-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
450	6,7-diF-Q	-CH ₂ O-	H	9-CH=CH ₂	-SCH ₂ CH ₂ COOH

451	6,7-diF-Q	-CH ₂ O-	H	8-CH=CH ₂	-SCH ₂ CH ₂ COOH
452	6,7-diF-Q	-CH ₂ O-	H	7-CH=CH ₂	-SCH ₂ CH ₂ COOH
453	6,7-diF-Q	-CH ₂ O-	H	9-NO ₂	-OCH ₂ COOH
454	6,7-diF-Q	-CH ₂ O-	H	9-NO ₂	-SCH ₂ CH ₂ COOH
455	6,7-diF-Q	-CH ₂ O-	H	8-NO ₂	-OCH ₂ COOH
456	6,7-diF-Q	-CH ₂ O-	H	8-NO ₂	-SCH ₂ CH ₂ COOH
457	6,7-diF-Q	-CH ₂ O-	H	7-NO ₂	-OCH ₂ COOH
458	6,7-diF-Q	-CH ₂ O-	H	7-NO ₂	-SCH ₂ CH ₂ COOH
459	6,7-diF-Q	-CH ₂ O-	1-F	H	-SCH ₂ CH ₂ COOH
460	6,7-diF-Q	-CH ₂ O-	2-F	H	-SCH ₂ CH ₂ COOH
461	6,7-diF-Q	-CH ₂ O-	1-Cl	H	-SCH ₂ CH ₂ COOH
462	6,7-diF-Q	-CH ₂ O-	2-Cl	H	-SCH ₂ CH ₂ COOH
463	6,7-diF-Q	-CH ₂ O-	1-CH ₃	H	-SCH ₂ CH ₂ COOH
464	6,7-diF-Q	-CH ₂ O-	2-CH ₃	H	-SCH ₂ CH ₂ COOH
465	6,7-diF-Q	-CH ₂ O-	1-OCH ₃	H	-SCH ₂ CH ₂ COOH
466	6,7-diF-Q	-CH ₂ O-	2-OCH ₃	H	-SCH ₂ CH ₂ COOH
467	6,7-diF-Q	-CH ₂ O-	1-NO ₂	H	-SCH ₂ CH ₂ COOH
468	6,7-diF-Q	-CH ₂ O-	2-NO ₂	H	-SCH ₂ CH ₂ COOH
469	6,7-diF-Q	-CH ₂ O-	1-CN	H	-SCH ₂ CH ₂ COOH
470	6,7-diF-Q	-CH ₂ O-	2-CN	H	-SCH ₂ CH ₂ COOH
471	7-Cl,6-F-Q	-CH=CH-	H	H	-OCH ₂ COOH
472	7-Cl,6-F-Q	-CH=CH-	H	H	-OCH(CH ₃)COOH
473	7-Cl,6-F-Q	-CH=CH-	H	H	-OCH ₂ CH ₂ COOH
474	7-Cl,6-F-Q	-CH=CH-	H	H	-OCH ₂ CH(CH ₃)COOH
475	7-Cl,6-F-Q	-CH=CH-	H	H	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
476	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ COOH
477	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH(CH ₃)COOH
478	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
479	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)COOH
480	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₃) ₂ COOH
481	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₂ CH ₃)COOH
482	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH(CH ₃)CH ₂ COOH
483	7-Cl,6-F-Q	-CH=CH-	H	H	-SC(CH ₃) ₂ CH ₂ COOH

484	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CH ₂ COOH
485	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)CH ₂ COOH
486	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₃) ₂ CH ₂ COOH
487	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
488	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CF ₂ COOH
489	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CF ₂ CH ₂ COOH
490	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ -Tet
491	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ -Tet
492	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ NHSO ₂ CF ₃
493	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ CH ₃
494	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ CF ₃
495	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ Ph
496	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ (2-CH ₃ -Ph)
497	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ NHSO ₂ CF ₃
498	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CH ₃
499	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CF ₃
500	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ Ph
501	7-Cl,6-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ (2-CH ₃ -Ph)
502	7-Cl,6-F-Q	-CH=CH-	H	H	-CH ₂ COOH
503	7-Cl,6-F-Q	-CH=CH-	H	H	-CH ₂ CH ₂ COOH
504	7-Cl,6-F-Q	-CH=CH-	H	H	-CH ₂ CH ₂ CH ₂ COOH
505	7-Cl,6-F-Q	-CH=CH-	H	9-F	-OCH ₂ COOH
506	7-Cl,6-F-Q	-CH=CH-	H	9-F	-OCH(CH ₃)COOH
507	7-Cl,6-F-Q	-CH=CH-	H	9-F	-OCH ₂ CH ₂ COOH
508	7-Cl,6-F-Q	-CH=CH-	H	9-F	-SCH ₂ COOH
509	7-Cl,6-F-Q	-CH=CH-	H	9-F	-SCH(CH ₃)COOH
510	7-Cl,6-F-Q	-CH=CH-	H	9-F	-SCH ₂ CH ₂ COOH
511	7-Cl,6-F-Q	-CH=CH-	H	9-F	-SCH ₂ CH(CH ₃)COOH
512	7-Cl,6-F-Q	-CH=CH-	H	9-F	-SCH ₂ CH(CH ₂ CH ₃)COOH
513	7-Cl,6-F-Q	-CH=CH-	H	9-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
514	7-Cl,6-F-Q	-CH=CH-	H	8-F	-OCH ₂ COOH
515	7-Cl,6-F-Q	-CH=CH-	H	8-F	-OCH ₂ CH ₂ COOH
516	7-Cl,6-F-Q	-CH=CH-	H	8-F	-SCH ₂ COOH

517	7-Cl,6-F-Q	-CH=CH-	H	8-F	-SCH ₂ CH ₂ COOH
518	7-Cl,6-F-Q	-CH=CH-	H	8-F	-SCH ₂ CH(CH ₃)COOH
519	7-Cl,6-F-Q	-CH=CH-	H	8-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
520	7-Cl,6-F-Q	-CH=CH-	H	7-F	-OCH ₂ COOH
521	7-Cl,6-F-Q	-CH=CH-	H	7-F	-OCH ₂ CH ₂ COOH
522	7-Cl,6-F-Q	-CH=CH-	H	7-F	-SCH ₂ COOH
523	7-Cl,6-F-Q	-CH=CH-	H	7-F	-SCH ₂ CH ₂ COOH
524	7-Cl,6-F-Q	-CH=CH-	H	7-F	-SCH ₂ CH(CH ₃)COOH
525	7-Cl,6-F-Q	-CH=CH-	H	6-F	-OCH ₂ COOH
526	7-Cl,6-F-Q	-CH=CH-	H	6-F	-OCH(CH ₃)COOH
527	7-Cl,6-F-Q	-CH=CH-	H	6-F	-OCH ₂ CH ₂ COOH
528	7-Cl,6-F-Q	-CH=CH-	H	6-F	-OCH ₂ CH(CH ₃)COOH
529	7-Cl,6-F-Q	-CH=CH-	H	6-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
530	7-Cl,6-F-Q	-CH=CH-	H	6-F	-SCH ₂ COOH
531	7-Cl,6-F-Q	-CH=CH-	H	6-F	-SCH(CH ₃)COOH
532	7-Cl,6-F-Q	-CH=CH-	H	6-F	-SCH ₂ CH ₂ COOH
533	7-Cl,6-F-Q	-CH=CH-	H	6-F	-SCH ₂ CH(CH ₃)COOH
534	7-Cl,6-F-Q	-CH=CH-	H	6-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
535	7-Cl,6-F-Q	-CH=CH-	H	9-Cl	-OCH ₂ COOH
536	7-Cl,6-F-Q	-CH=CH-	H	9-Cl	-OCH(CH ₃)COOH
537	7-Cl,6-F-Q	-CH=CH-	H	9-Cl	-OCH ₂ CH ₂ COOH
538	7-Cl,6-F-Q	-CH=CH-	H	9-Cl	-OCH ₂ CH(CH ₃)COOH
539	7-Cl,6-F-Q	-CH=CH-	H	9-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
540	7-Cl,6-F-Q	-CH=CH-	H	9-Cl	-SCH ₂ COOH
541	7-Cl,6-F-Q	-CH=CH-	H	9-Cl	-SCH(CH ₃)COOH
542	7-Cl,6-F-Q	-CH=CH-	H	9-Cl	-SCH ₂ CH ₂ COOH
543	7-Cl,6-F-Q	-CH=CH-	H	9-Cl	-SCH ₂ CH(CH ₃)COOH
544	7-Cl,6-F-Q	-CH=CH-	H	9-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
545	7-Cl,6-F-Q	-CH=CH-	H	8-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
546	7-Cl,6-F-Q	-CH=CH-	H	8-Cl	-OCH ₂ COOH
547	7-Cl,6-F-Q	-CH=CH-	H	8-Cl	-OCH ₂ CH ₂ COOH
548	7-Cl,6-F-Q	-CH=CH-	H	8-Cl	-SCH ₂ COOH
549	7-Cl,6-F-Q	-CH=CH-	H	8-Cl	-SCH ₂ CH ₂ COOH

550	7-Cl,6-F-Q	-CH=CH-	H	8-Cl	-SCH ₂ CH(CH ₃)COOH
551	7-Cl,6-F-Q	-CH=CH-	H	7-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
552	7-Cl,6-F-Q	-CH=CH-	H	7-Cl	-OCH ₂ COOH
553	7-Cl,6-F-Q	-CH=CH-	H	7-Cl	-OCH ₂ CH ₂ COOH
554	7-Cl,6-F-Q	-CH=CH-	H	7-Cl	-SCH ₂ COOH
555	7-Cl,6-F-Q	-CH=CH-	H	7-Cl	-SCH ₂ CH ₂ COOH
556	7-Cl,6-F-Q	-CH=CH-	H	7-Cl	-SCH ₂ CH(CH ₃)COOH
557	7-Cl,6-F-Q	-CH=CH-	H	6-Cl	-OCH ₂ COOH
558	7-Cl,6-F-Q	-CH=CH-	H	6-Cl	-OCH(CH ₃)COOH
559	7-Cl,6-F-Q	-CH=CH-	H	6-Cl	-OCH ₂ CH ₂ COOH
560	7-Cl,6-F-Q	-CH=CH-	H	6-Cl	-OCH ₂ CH(CH ₃)COOH
561	7-Cl,6-F-Q	-CH=CH-	H	6-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
562	7-Cl,6-F-Q	-CH=CH-	H	6-Cl	-SCH ₂ COOH
563	7-Cl,6-F-Q	-CH=CH-	H	6-Cl	-SCH(CH ₃)COOH
564	7-Cl,6-F-Q	-CH=CH-	H	6-Cl	-SCH ₂ CH ₂ COOH
565	7-Cl,6-F-Q	-CH=CH-	H	6-Cl	-SCH ₂ CH(CH ₃)COOH
566	7-Cl,6-F-Q	-CH=CH-	H	6-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
567	7-Cl,6-F-Q	-CH=CH-	H	9-CN	-OCH ₂ COOH
568	7-Cl,6-F-Q	-CH=CH-	H	9-CN	-OCH(CH ₃)COOH
569	7-Cl,6-F-Q	-CH=CH-	H	9-CN	-OCH ₂ CH ₂ COOH
570	7-Cl,6-F-Q	-CH=CH-	H	9-CN	-OCH ₂ CH(CH ₃)COOH
571	7-Cl,6-F-Q	-CH=CH-	H	9-CN	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
572	7-Cl,6-F-Q	-CH=CH-	H	9-CN	-SCH ₂ COOH
573	7-Cl,6-F-Q	-CH=CH-	H	9-CN	-SCH(CH ₃)COOH
574	7-Cl,6-F-Q	-CH=CH-	H	9-CN	-SCH ₂ CH ₂ COOH
575	7-Cl,6-F-Q	-CH=CH-	H	9-CN	-SCH ₂ CH(CH ₃)COOH
576	7-Cl,6-F-Q	-CH=CH-	H	9-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
577	7-Cl,6-F-Q	-CH=CH-	H	8-CN	-OCH ₂ COOH
578	7-Cl,6-F-Q	-CH=CH-	H	8-CN	-OCH ₂ CH ₂ COOH
579	7-Cl,6-F-Q	-CH=CH-	H	8-CN	-SCH ₂ COOH
580	7-Cl,6-F-Q	-CH=CH-	H	8-CN	-SCH ₂ CH ₂ COOH
581	7-Cl,6-F-Q	-CH=CH-	H	8-CN	-SCH ₂ CH(CH ₃)COOH
582	7-Cl,6-F-Q	-CH=CH-	H	8-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH

583	7-Cl,6-F-Q	-CH=CH-	H	7-CN	-OCH ₂ COOH
584	7-Cl,6-F-Q	-CH=CH-	H	7-CN	-OCH ₂ CH ₂ COOH
585	7-Cl,6-F-Q	-CH=CH-	H	7-CN	-SCH ₂ COOH
586	7-Cl,6-F-Q	-CH=CH-	H	7-CN	-SCH ₂ CH ₂ COOH
587	7-Cl,6-F-Q	-CH=CH-	H	7-CN	-SCH ₂ CH(CH ₃)COOH
588	7-Cl,6-F-Q	-CH=CH-	H	9-CH ₃	-OCH ₂ COOH
589	7-Cl,6-F-Q	-CH=CH-	H	9-CH ₃	-OCH ₂ CH ₂ COOH
590	7-Cl,6-F-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ COOH
591	7-Cl,6-F-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH ₂ COOH
592	7-Cl,6-F-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH(CH ₃)COOH
593	7-Cl,6-F-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH(CH ₂ CH ₃)COOH
594	7-Cl,6-F-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
595	7-Cl,6-F-Q	-CH=CH-	H	8-CH ₃	-OCH ₂ COOH
596	7-Cl,6-F-Q	-CH=CH-	H	8-CH ₃	-OCH ₂ CH ₂ COOH
597	7-Cl,6-F-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ COOH
598	7-Cl,6-F-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ CH ₂ COOH
599	7-Cl,6-F-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ CH(CH ₃)COOH
600	7-Cl,6-F-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
601	7-Cl,6-F-Q	-CH=CH-	H	7-CH ₃	-OCH ₂ COOH
602	7-Cl,6-F-Q	-CH=CH-	H	7-CH ₃	-OCH ₂ CH ₂ COOH
603	7-Cl,6-F-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ COOH
604	7-Cl,6-F-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ CH ₂ COOH
605	7-Cl,6-F-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ CH(CH ₃)COOH
606	7-Cl,6-F-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ COOH
607	7-Cl,6-F-Q	-CH=CH-	H	9-CF ₃	-OCH(CH ₃)COOH
608	7-Cl,6-F-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ CH ₂ COOH
609	7-Cl,6-F-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ CH(CH ₃)COOH
610	7-Cl,6-F-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
611	7-Cl,6-F-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ COOH
612	7-Cl,6-F-Q	-CH=CH-	H	9-CF ₃	-SCH(CH ₃)COOH
613	7-Cl,6-F-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ CH ₂ COOH
614	7-Cl,6-F-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ CH(CH ₃)COOH
615	7-Cl,6-F-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH

616	7-Cl,6-F-Q	-CH=CH-	H	8-CF ₃	-OCH ₂ COOH
617	7-Cl,6-F-Q	-CH=CH-	H	8-CF ₃	-OCH ₂ CH ₂ COOH
618	7-Cl,6-F-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ COOH
619	7-Cl,6-F-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ CH ₂ COOH
620	7-Cl,6-F-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ CH(CH ₃)COOH
621	7-Cl,6-F-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
622	7-Cl,6-F-Q	-CH=CH-	H	7-CF ₃	-OCH ₂ COOH
623	7-Cl,6-F-Q	-CH=CH-	H	7-CF ₃	-OCH ₂ CH ₂ COOH
624	7-Cl,6-F-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ COOH
625	7-Cl,6-F-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ CH ₂ COOH
626	7-Cl,6-F-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ CH(CH ₃)COOH
627	7-Cl,6-F-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ COOH
628	7-Cl,6-F-Q	-CH=CH-	H	9-C≡CH	-OCH(CH ₃)COOH
629	7-Cl,6-F-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ CH ₂ COOH
630	7-Cl,6-F-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ CH(CH ₃)COOH
631	7-Cl,6-F-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
632	7-Cl,6-F-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ COOH
633	7-Cl,6-F-Q	-CH=CH-	H	9-C≡CH	-SCH(CH ₃)COOH
634	7-Cl,6-F-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ CH ₂ COOH
635	7-Cl,6-F-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ CH(CH ₃)COOH
636	7-Cl,6-F-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
637	7-Cl,6-F-Q	-CH=CH-	H	8-C≡CH	-OCH ₂ COOH
638	7-Cl,6-F-Q	-CH=CH-	H	8-C≡CH	-OCH ₂ CH ₂ COOH
639	7-Cl,6-F-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ COOH
640	7-Cl,6-F-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ CH ₂ COOH
641	7-Cl,6-F-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ CH(CH ₃)COOH
642	7-Cl,6-F-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
643	7-Cl,6-F-Q	-CH=CH-	H	7-C≡CH	-OCH ₂ COOH
644	7-Cl,6-F-Q	-CH=CH-	H	7-C≡CH	-OCH ₂ CH ₂ COOH
645	7-Cl,6-F-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ COOH
646	7-Cl,6-F-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ CH ₂ COOH
647	7-Cl,6-F-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ CH(CH ₃)COOH
648	7-Cl,6-F-Q	-CH=CH-	H	9-CH ₂ OH	-OCH ₂ COOH

649	7-Cl,6-F-Q	-CH=CH-	H	9-CH ₂ OH	-SCH ₂ CH ₂ COOH
650	7-Cl,6-F-Q	-CH=CH-	H	8-CH ₂ OH	-OCH ₂ COOH
651	7-Cl,6-F-Q	-CH=CH-	H	8-CH ₂ OH	-SCH ₂ CH ₂ COOH
652	7-Cl,6-F-Q	-CH=CH-	H	7-CH ₂ OH	-OCH ₂ COOH
653	7-Cl,6-F-Q	-CH=CH-	H	7-CH ₂ OH	-SCH ₂ CH ₂ COOH
654	7-Cl,6-F-Q	-CH=CH-	H	9- C(CH ₃) ₂ OH	-OCH ₂ COOH
655	7-Cl,6-F-Q	-CH=CH-	H	9- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
656	7-Cl,6-F-Q	-CH=CH-	H	8- C(CH ₃) ₂ OH	-OCH ₂ COOH
657	7-Cl,6-F-Q	-CH=CH-	H	8- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
658	7-Cl,6-F-Q	-CH=CH-	H	7- C(CH ₃) ₂ OH	-OCH ₂ COOH
659	7-Cl,6-F-Q	-CH=CH-	H	7- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
660	7-Cl,6-F-Q	-CH=CH-	H	9-OCH ₃	-OCH ₂ COOH
661	7-Cl,6-F-Q	-CH=CH-	H	9-OCH ₃	-SCH ₂ CH ₂ COOH
662	7-Cl,6-F-Q	-CH=CH-	H	8-OCH ₃	-OCH ₂ COOH
663	7-Cl,6-F-Q	-CH=CH-	H	8-OCH ₃	-SCH ₂ CH ₂ COOH
664	7-Cl,6-F-Q	-CH=CH-	H	7-OCH ₃	-OCH ₂ COOH
665	7-Cl,6-F-Q	-CH=CH-	H	7-OCH ₃	-SCH ₂ CH ₂ COOH
666	7-Cl,6-F-Q	-CH=CH-	H	9-OCHF ₂	-OCH ₂ COOH
667	7-Cl,6-F-Q	-CH=CH-	H	9-OCHF ₂	-SCH ₂ CH ₂ COOH
668	7-Cl,6-F-Q	-CH=CH-	H	8-OCHF ₂	-OCH ₂ COOH
669	7-Cl,6-F-Q	-CH=CH-	H	8-OCHF ₂	-SCH ₂ CH ₂ COOH
670	7-Cl,6-F-Q	-CH=CH-	H	7-OCHF ₂	-OCH ₂ COOH
671	7-Cl,6-F-Q	-CH=CH-	H	7-OCHF ₂	-SCH ₂ CH ₂ COOH
672	7-Cl,6-F-Q	-CH=CH-	H	9-OCF ₃	-OCH ₂ COOH
673	7-Cl,6-F-Q	-CH=CH-	H	9-OCF ₃	-SCH ₂ CH ₂ COOH
674	7-Cl,6-F-Q	-CH=CH-	H	8-OCF ₃	-OCH ₂ COOH
675	7-Cl,6-F-Q	-CH=CH-	H	8-OCF ₃	-SCH ₂ CH ₂ COOH

676	7-Cl,6-F-Q	-CH=CH-	H	7-OCF ₃	-OCH ₂ COOH
677	7-Cl,6-F-Q	-CH=CH-	H	7-OCF ₃	-SCH ₂ CH ₂ COOH
678	7-Cl,6-F-Q	-CH=CH-	H	9-SOCH ₃	-OCH ₂ COOH
679	7-Cl,6-F-Q	-CH=CH-	H	9-SOCH ₃	-SCH ₂ CH ₂ COOH
680	7-Cl,6-F-Q	-CH=CH-	H	8-SOCH ₃	-OCH ₂ COOH
681	7-Cl,6-F-Q	-CH=CH-	H	8-SOCH ₃	-SCH ₂ CH ₂ COOH
682	7-Cl,6-F-Q	-CH=CH-	H	7-SOCH ₃	-OCH ₂ COOH
683	7-Cl,6-F-Q	-CH=CH-	H	7-SOCH ₃	-SCH ₂ CH ₂ COOH
684	7-Cl,6-F-Q	-CH=CH-	H	9-SO ₂ CH ₃	-OCH ₂ COOH
685	7-Cl,6-F-Q	-CH=CH-	H	9-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
686	7-Cl,6-F-Q	-CH=CH-	H	8-SO ₂ CH ₃	-OCH ₂ COOH
687	7-Cl,6-F-Q	-CH=CH-	H	8-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
688	7-Cl,6-F-Q	-CH=CH-	H	7-SO ₂ CH ₃	-OCH ₂ COOH
689	7-Cl,6-F-Q	-CH=CH-	H	7-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
690	7-Cl,6-F-Q	-CH=CH-	H	9-CH=CH ₂	-SCH ₂ CH ₂ COOH
691	7-Cl,6-F-Q	-CH=CH-	H	8-CH=CH ₂	-SCH ₂ CH ₂ COOH
692	7-Cl,6-F-Q	-CH=CH-	H	7-CH=CH ₂	-SCH ₂ CH ₂ COOH
693	7-Cl,6-F-Q	-CH=CH-	H	9-NO ₂	-OCH ₂ COOH
694	7-Cl,6-F-Q	-CH=CH-	H	9-NO ₂	-SCH ₂ CH ₂ COOH
695	7-Cl,6-F-Q	-CH=CH-	H	8-NO ₂	-OCH ₂ COOH
696	7-Cl,6-F-Q	-CH=CH-	H	8-NO ₂	-SCH ₂ CH ₂ COOH
697	7-Cl,6-F-Q	-CH=CH-	H	7-NO ₂	-OCH ₂ COOH
698	7-Cl,6-F-Q	-CH=CH-	H	7-NO ₂	-SCH ₂ CH ₂ COOH
699	7-Cl,6-F-Q	-CH=CH-	1-F	H	-SCH ₂ CH ₂ COOH
700	7-Cl,6-F-Q	-CH=CH-	2-F	H	-SCH ₂ CH ₂ COOH
701	7-Cl,6-F-Q	-CH=CH-	1-Cl	H	-SCH ₂ CH ₂ COOH
702	7-Cl,6-F-Q	-CH=CH-	2-Cl	H	-SCH ₂ CH ₂ COOH
703	7-Cl,6-F-Q	-CH=CH-	1-CH ₃	H	-SCH ₂ CH ₂ COOH
704	7-Cl,6-F-Q	-CH=CH-	2-CH ₃	H	-SCH ₂ CH ₂ COOH
705	7-Cl,6-F-Q	-CH=CH-	1-OCH ₃	H	-SCH ₂ CH ₂ COOH
706	7-Cl,6-F-Q	-CH=CH-	2-OCH ₃	H	-SCH ₂ CH ₂ COOH
707	7-Cl,6-F-Q	-CH=CH-	1-NO ₂	H	-SCH ₂ CH ₂ COOH
708	7-Cl,6-F-Q	-CH=CH-	2-NO ₂	H	-SCH ₂ CH ₂ COOH

709	7-Cl,6-F-Q	-CH=CH-	1-CN	H	-SCH ₂ CH ₂ COOH
710	7-Cl,6-F-Q	-CH=CH-	2-CN	H	-SCH ₂ CH ₂ COOH
711	7-Cl,6-F-Q	-CH ₂ O-	H	H	-OCH ₂ COOH
712	7-Cl,6-F-Q	-CH ₂ O-	H	H	-OCH(CH ₃)COOH
713	7-Cl,6-F-Q	-CH ₂ O-	H	H	-OCH ₂ CH ₂ COOH
714	7-Cl,6-F-Q	-CH ₂ O-	H	H	-OCH ₂ CH(CH ₃)COOH
715	7-Cl,6-F-Q	-CH ₂ O-	H	H	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
716	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ COOH
717	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH(CH ₃)COOH
718	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
719	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)COOH
720	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₃) ₂ COOH
721	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₂ CH ₃)COOH
722	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH(CH ₃)CH ₂ COOH
723	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SC(CH ₃) ₂ CH ₂ COOH
724	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CH ₂ COOH
725	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)CH ₂ COOH
726	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₃) ₂ CH ₂ COOH
727	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
728	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CF ₂ COOH
729	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CF ₂ CH ₂ COOH
730	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ -Tet
731	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ -Tet
732	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ NHSO ₂ CF ₃
733	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ CH ₃
734	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ CF ₃
735	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ Ph
736	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ (2-CH ₃ -Ph)
737	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ NHSO ₂ CF ₃
738	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CH ₃
739	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CF ₃
740	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ Ph
741	7-Cl,6-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ (2-CH ₃ -Ph)

742	7-Cl,6-F-Q	-CH ₂ O-	H	H	-CH ₂ COOH
743	7-Cl,6-F-Q	-CH ₂ O-	H	H	-CH ₂ CH ₂ COOH
744	7-Cl,6-F-Q	-CH ₂ O-	H	H	-CH ₂ CH ₂ CH ₂ COOH
745	7-Cl,6-F-Q	-CH ₂ O-	H	9-F	-OCH ₂ COOH
746	7-Cl,6-F-Q	-CH ₂ O-	H	9-F	-OCH(CH ₃)COOH
747	7-Cl,6-F-Q	-CH ₂ O-	H	9-F	-OCH ₂ CH ₂ COOH
748	7-Cl,6-F-Q	-CH ₂ O-	H	9-F	-OCH ₂ CH(CH ₃)COOH
749	7-Cl,6-F-Q	-CH ₂ O-	H	9-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
750	7-Cl,6-F-Q	-CH ₂ O-	H	9-F	-SCH ₂ COOH
751	7-Cl,6-F-Q	-CH ₂ O-	H	9-F	-SCH(CH ₃)COOH
752	7-Cl,6-F-Q	-CH ₂ O-	H	9-F	-SCH ₂ CH ₂ COOH
753	7-Cl,6-F-Q	-CH ₂ O-	H	9-F	-SCH ₂ CH(CH ₃)COOH
754	7-Cl,6-F-Q	-CH ₂ O-	H	9-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
755	7-Cl,6-F-Q	-CH ₂ O-	H	8-F	-OCH ₂ COOH
756	7-Cl,6-F-Q	-CH ₂ O-	H	8-F	-OCH ₂ CH ₂ COOH
757	7-Cl,6-F-Q	-CH ₂ O-	H	8-F	-SCH ₂ COOH
758	7-Cl,6-F-Q	-CH ₂ O-	H	8-F	-SCH ₂ CH ₂ COOH
759	7-Cl,6-F-Q	-CH ₂ O-	H	8-F	-SCH ₂ CH(CH ₃)COOH
760	7-Cl,6-F-Q	-CH ₂ O-	H	8-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
761	7-Cl,6-F-Q	-CH ₂ O-	H	7-F	-OCH ₂ COOH
762	7-Cl,6-F-Q	-CH ₂ O-	H	7-F	-OCH ₂ CH ₂ COOH
763	7-Cl,6-F-Q	-CH ₂ O-	H	7-F	-SCH ₂ COOH
764	7-Cl,6-F-Q	-CH ₂ O-	H	7-F	-SCH ₂ CH ₂ COOH
765	7-Cl,6-F-Q	-CH ₂ O-	H	7-F	-SCH ₂ CH(CH ₃)COOH
766	7-Cl,6-F-Q	-CH ₂ O-	H	6-F	-OCH ₂ COOH
767	7-Cl,6-F-Q	-CH ₂ O-	H	6-F	-OCH(CH ₃)COOH
768	7-Cl,6-F-Q	-CH ₂ O-	H	6-F	-OCH ₂ CH ₂ COOH
769	7-Cl,6-F-Q	-CH ₂ O-	H	6-F	-OCH ₂ CH(CH ₃)COOH
770	7-Cl,6-F-Q	-CH ₂ O-	H	6-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
771	7-Cl,6-F-Q	-CH ₂ O-	H	6-F	-SCH ₂ COOH
772	7-Cl,6-F-Q	-CH ₂ O-	H	6-F	-SCH(CH ₃)COOH
773	7-Cl,6-F-Q	-CH ₂ O-	H	6-F	-SCH ₂ CH ₂ COOH
774	7-Cl,6-F-Q	-CH ₂ O-	H	6-F	-SCH ₂ CH(CH ₃)COOH

775	7-Cl,6-F-Q	-CH ₂ O-	H	6-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
776	7-Cl,6-F-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ COOH
777	7-Cl,6-F-Q	-CH ₂ O-	H	9-Cl	-OCH(CH ₃)COOH
778	7-Cl,6-F-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ CH ₂ COOH
779	7-Cl,6-F-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ CH(CH ₃)COOH
780	7-Cl,6-F-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
781	7-Cl,6-F-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ COOH
782	7-Cl,6-F-Q	-CH ₂ O-	H	9-Cl	-SCH(CH ₃)COOH
783	7-Cl,6-F-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ CH ₂ COOH
784	7-Cl,6-F-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ CH(CH ₃)COOH
785	7-Cl,6-F-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
786	7-Cl,6-F-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
787	7-Cl,6-F-Q	-CH ₂ O-	H	8-Cl	-OCH ₂ COOH
788	7-Cl,6-F-Q	-CH ₂ O-	H	8-Cl	-OCH ₂ CH ₂ COOH
789	7-Cl,6-F-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ COOH
790	7-Cl,6-F-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ CH ₂ COOH
791	7-Cl,6-F-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ CH(CH ₃)COOH
792	7-Cl,6-F-Q	-CH ₂ O-	H	7-Cl	-OCH ₂ COOH
793	7-Cl,6-F-Q	-CH ₂ O-	H	7-Cl	-OCH ₂ CH ₂ COOH
794	7-Cl,6-F-Q	-CH ₂ O-	H	7-Cl	-SCH ₂ COOH
795	7-Cl,6-F-Q	-CH ₂ O-	H	7-Cl	-SCH ₂ CH ₂ COOH
796	7-Cl,6-F-Q	-CH ₂ O-	H	7-Cl	-SCH ₂ CH(CH ₃)COOH
797	7-Cl,6-F-Q	-CH ₂ O-	H	9-CN	-OCH ₂ COOH
798	7-Cl,6-F-Q	-CH ₂ O-	H	9-CN	-OCH(CH ₃)COOH
799	7-Cl,6-F-Q	-CH ₂ O-	H	9-CN	-OCH ₂ CH ₂ COOH
800	7-Cl,6-F-Q	-CH ₂ O-	H	9-CN	-OCH ₂ CH(CH ₃)COOH
801	7-Cl,6-F-Q	-CH ₂ O-	H	9-CN	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
802	7-Cl,6-F-Q	-CH ₂ O-	H	9-CN	-SCH ₂ COOH
803	7-Cl,6-F-Q	-CH ₂ O-	H	9-CN	-SCH(CH ₃)COOH
804	7-Cl,6-F-Q	-CH ₂ O-	H	9-CN	-SCH ₂ CH ₂ COOH
805	7-Cl,6-F-Q	-CH ₂ O-	H	9-CN	-SCH ₂ CH(CH ₃)COOH
806	7-Cl,6-F-Q	-CH ₂ O-	H	9-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
807	7-Cl,6-F-Q	-CH ₂ O-	H	8-CN	-OCH ₂ COOH

808	7-Cl,6-F-Q	-CH ₂ O-	H	8-CN	-OCH ₂ CH ₂ COOH
809	7-Cl,6-F-Q	-CH ₂ O-	H	8-CN	-SCH ₂ COOH
810	7-Cl,6-F-Q	-CH ₂ O-	H	8-CN	-SCH ₂ CH ₂ COOH
811	7-Cl,6-F-Q	-CH ₂ O-	H	8-CN	-SCH ₂ CH(CH ₃)COOH
812	7-Cl,6-F-Q	-CH ₂ O-	H	8-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
813	7-Cl,6-F-Q	-CH ₂ O-	H	7-CN	-OCH ₂ COOH
814	7-Cl,6-F-Q	-CH ₂ O-	H	7-CN	-OCH ₂ CH ₂ COOH
815	7-Cl,6-F-Q	-CH ₂ O-	H	7-CN	-SCH ₂ COOH
816	7-Cl,6-F-Q	-CH ₂ O-	H	7-CN	-SCH ₂ CH ₂ COOH
817	7-Cl,6-F-Q	-CH ₂ O-	H	7-CN	-SCH ₂ CH(CH ₃)COOH
818	7-Cl,6-F-Q	-CH ₂ O-	H	9-CH ₃	-OCH ₂ COOH
819	7-Cl,6-F-Q	-CH ₂ O-	H	9-CH ₃	-OCH ₂ CH ₂ COOH
820	7-Cl,6-F-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ COOH
821	7-Cl,6-F-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ CH ₂ COOH
822	7-Cl,6-F-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ CH(CH ₃)COOH
823	7-Cl,6-F-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ CH(CH ₂ CH ₃)COOH
824	7-Cl,6-F-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
825	7-Cl,6-F-Q	-CH ₂ O-	H	8-CH ₃	-OCH ₂ COOH
826	7-Cl,6-F-Q	-CH ₂ O-	H	8-CH ₃	-OCH ₂ CH ₂ COOH
827	7-Cl,6-F-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ COOH
828	7-Cl,6-F-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ CH ₂ COOH
829	7-Cl,6-F-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ CH(CH ₃)COOH
830	7-Cl,6-F-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
831	7-Cl,6-F-Q	-CH ₂ O-	H	7-CH ₃	-OCH ₂ COOH
832	7-Cl,6-F-Q	-CH ₂ O-	H	7-CH ₃	-OCH ₂ CH ₂ COOH
833	7-Cl,6-F-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ COOH
834	7-Cl,6-F-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ CH ₂ COOH
835	7-Cl,6-F-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ CH(CH ₃)COOH
836	7-Cl,6-F-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ COOH
837	7-Cl,6-F-Q	-CH ₂ O-	H	9-CF ₃	-OCH(CH ₃)COOH
838	7-Cl,6-F-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ CH ₂ COOH
839	7-Cl,6-F-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ CH(CH ₃)COOH
840	7-Cl,6-F-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH

841	7-Cl,6-F-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ COOH
842	7-Cl,6-F-Q	-CH ₂ O-	H	9-CF ₃	-SCH(CH ₃)COOH
843	7-Cl,6-F-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ CH ₂ COOH
844	7-Cl,6-F-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ CH(CH ₃)COOH
845	7-Cl,6-F-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
846	7-Cl,6-F-Q	-CH ₂ O-	H	8-CF ₃	-OCH ₂ COOH
847	7-Cl,6-F-Q	-CH ₂ O-	H	8-CF ₃	-OCH ₂ CH ₂ COOH
848	7-Cl,6-F-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ COOH
849	7-Cl,6-F-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ CH ₂ COOH
850	7-Cl,6-F-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ CH(CH ₃)COOH
851	7-Cl,6-F-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
852	7-Cl,6-F-Q	-CH ₂ O-	H	7-CF ₃	-OCH ₂ COOH
853	7-Cl,6-F-Q	-CH ₂ O-	H	7-CF ₃	-OCH ₂ CH ₂ COOH
854	7-Cl,6-F-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ COOH
855	7-Cl,6-F-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ CH ₂ COOH
856	7-Cl,6-F-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ CH(CH ₃)COOH
857	7-Cl,6-F-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ COOH
858	7-Cl,6-F-Q	-CH ₂ O-	H	9-C≡CH	-OCH(CH ₃)COOH
859	7-Cl,6-F-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ CH ₂ COOH
860	7-Cl,6-F-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ CH(CH ₃)COOH
861	7-Cl,6-F-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
862	7-Cl,6-F-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ COOH
863	7-Cl,6-F-Q	-CH ₂ O-	H	9-C≡CH	-SCH(CH ₃)COOH
864	7-Cl,6-F-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ CH ₂ COOH
865	7-Cl,6-F-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ CH(CH ₃)COOH
866	7-Cl,6-F-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
867	7-Cl,6-F-Q	-CH ₂ O-	H	8-C≡CH	-OCH ₂ COOH
868	7-Cl,6-F-Q	-CH ₂ O-	H	8-C≡CH	-OCH ₂ CH ₂ COOH
869	7-Cl,6-F-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ COOH
870	7-Cl,6-F-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ CH ₂ COOH
871	7-Cl,6-F-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ CH(CH ₃)COOH
872	7-Cl,6-F-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
873	7-Cl,6-F-Q	-CH ₂ O-	H	7-C≡CH	-OCH ₂ COOH

874	7-Cl,6-F-Q	-CH ₂ O-	H	7-C≡CH	-OCH ₂ CH ₂ COOH
875	7-Cl,6-F-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ COOH
876	7-Cl,6-F-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ CH ₂ COOH
877	7-Cl,6-F-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ CH(CH ₃)COOH
878	7-Cl,6-F-Q	-CH ₂ O-	H	9-CH ₂ OH	-OCH ₂ COOH
879	7-Cl,6-F-Q	-CH ₂ O-	H	9-CH ₂ OH	-SCH ₂ CH ₂ COOH
880	7-Cl,6-F-Q	-CH ₂ O-	H	8-CH ₂ OH	-OCH ₂ COOH
881	7-Cl,6-F-Q	-CH ₂ O-	H	8-CH ₂ OH	-SCH ₂ CH ₂ COOH
885	7-Cl,6-F-Q	-CH ₂ O-	H	9- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
886	7-Cl,6-F-Q	-CH ₂ O-	H	8- C(CH ₃) ₂ OH	-OCH ₂ COOH
887	7-Cl,6-F-Q	-CH ₂ O-	H	8- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
888	7-Cl,6-F-Q	-CH ₂ O-	H	7- C(CH ₃) ₂ OH	-OCH ₂ COOH
889	7-Cl,6-F-Q	-CH ₂ O-	H	7- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
890	7-Cl,6-F-Q	-CH ₂ O-	H	9-OCH ₃	-OCH ₂ COOH
891	7-Cl,6-F-Q	-CH ₂ O-	H	9-OCH ₃	-SCH ₂ CH ₂ COOH
892	7-Cl,6-F-Q	-CH ₂ O-	H	8-OCH ₃	-OCH ₂ COOH
893	7-Cl,6-F-Q	-CH ₂ O-	H	8-OCH ₃	-SCH ₂ CH ₂ COOH
894	7-Cl,6-F-Q	-CH ₂ O-	H	7-OCH ₃	-OCH ₂ COOH
895	7-Cl,6-F-Q	-CH ₂ O-	H	7-OCH ₃	-SCH ₂ CH ₂ COOH
896	7-Cl,6-F-Q	-CH ₂ O-	H	9-OCHF ₂	-OCH ₂ COOH
897	7-Cl,6-F-Q	-CH ₂ O-	H	9-OCHF ₂	-SCH ₂ CH ₂ COOH
898	7-Cl,6-F-Q	-CH ₂ O-	H	8-OCHF ₂	-OCH ₂ COOH
899	7-Cl,6-F-Q	-CH ₂ O-	H	8-OCHF ₂	-SCH ₂ CH ₂ COOH
900	7-Cl,6-F-Q	-CH ₂ O-	H	7-OCHF ₂	-OCH ₂ COOH
901	7-Cl,6-F-Q	-CH ₂ O-	H	7-OCHF ₂	-SCH ₂ CH ₂ COOH
902	7-Cl,6-F-Q	-CH ₂ O-	H	9-OCF ₃	-OCH ₂ COOH
903	7-Cl,6-F-Q	-CH ₂ O-	H	9-OCF ₃	-SCH ₂ CH ₂ COOH
904	7-Cl,6-F-Q	-CH ₂ O-	H	8-OCF ₃	-OCH ₂ COOH

905	7-Cl,6-F-Q	-CH ₂ O-	H	8-OCF ₃	-SCH ₂ CH ₂ COOH
906	7-Cl,6-F-Q	-CH ₂ O-	H	7-OCF ₃	-OCH ₂ COOH
907	7-Cl,6-F-Q	-CH ₂ O-	H	7-OCF ₃	-SCH ₂ CH ₂ COOH
908	7-Cl,6-F-Q	-CH ₂ O-	H	9-SOCH ₃	-OCH ₂ COOH
909	7-Cl,6-F-Q	-CH ₂ O-	H	9-SOCH ₃	-SCH ₂ CH ₂ COOH
910	7-Cl,6-F-Q	-CH ₂ O-	H	8-SOCH ₃	-OCH ₂ COOH
911	7-Cl,6-F-Q	-CH ₂ O-	H	8-SOCH ₃	-SCH ₂ CH ₂ COOH
912	7-Cl,6-F-Q	-CH ₂ O-	H	7-SOCH ₃	-OCH ₂ COOH
913	7-Cl,6-F-Q	-CH ₂ O-	H	7-SOCH ₃	-SCH ₂ CH ₂ COOH
914	7-Cl,6-F-Q	-CH ₂ O-	H	9-SO ₂ CH ₃	-OCH ₂ COOH
915	7-Cl,6-F-Q	-CH ₂ O-	H	9-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
916	7-Cl,6-F-Q	-CH ₂ O-	H	8-SO ₂ CH ₃	-OCH ₂ COOH
917	7-Cl,6-F-Q	-CH ₂ O-	H	8-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
918	7-Cl,6-F-Q	-CH ₂ O-	H	7-SO ₂ CH ₃	-OCH ₂ COOH
919	7-Cl,6-F-Q	-CH ₂ O-	H	7-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
920	7-Cl,6-F-Q	-CH ₂ O-	H	9-CH=CH ₂	-SCH ₂ CH ₂ COOH
921	7-Cl,6-F-Q	-CH ₂ O-	H	8-CH=CH ₂	-SCH ₂ CH ₂ COOH
922	7-Cl,6-F-Q	-CH ₂ O-	H	7-CH=CH ₂	-SCH ₂ CH ₂ COOH
923	7-Cl,6-F-Q	-CH ₂ O-	H	9-NO ₂	-OCH ₂ COOH
924	7-Cl,6-F-Q	-CH ₂ O-	H	9-NO ₂	-SCH ₂ CH ₂ COOH
925	7-Cl,6-F-Q	-CH ₂ O-	H	8-NO ₂	-OCH ₂ COOH
926	7-Cl,6-F-Q	-CH ₂ O-	H	8-NO ₂	-SCH ₂ CH ₂ COOH
927	7-Cl,6-F-Q	-CH ₂ O-	H	7-NO ₂	-OCH ₂ COOH
928	7-Cl,6-F-Q	-CH ₂ O-	H	7-NO ₂	-SCH ₂ CH ₂ COOH
929	7-Cl,6-F-Q	-CH ₂ O-	1-F	H	-SCH ₂ CH ₂ COOH
930	7-Cl,6-F-Q	-CH ₂ O-	2-F	H	-SCH ₂ CH ₂ COOH
931	7-Cl,6-F-Q	-CH ₂ O-	1-Cl	H	-SCH ₂ CH ₂ COOH
932	7-Cl,6-F-Q	-CH ₂ O-	2-Cl	H	-SCH ₂ CH ₂ COOH
933	7-Cl,6-F-Q	-CH ₂ O-	1-CH ₃	H	-SCH ₂ CH ₂ COOH
934	7-Cl,6-F-Q	-CH ₂ O-	2-CH ₃	H	-SCH ₂ CH ₂ COOH
935	7-Cl,6-F-Q	-CH ₂ O-	1-OCH ₃	H	-SCH ₂ CH ₂ COOH
936	7-Cl,6-F-Q	-CH ₂ O-	2-OCH ₃	H	-SCH ₂ CH ₂ COOH
937	7-Cl,6-F-Q	-CH ₂ O-	1-NO ₂	H	-SCH ₂ CH ₂ COOH

938	7-Cl,6-F-Q	-CH ₂ O-	2-NO ₂	H	-SCH ₂ CH ₂ COOH
939	7-Cl,6-F-Q	-CH ₂ O-	1-CN	H	-SCH ₂ CH ₂ COOH
940	7-Cl,6-F-Q	-CH ₂ O-	2-CN	H	-SCH ₂ CH ₂ COOH
941	7-F-Q	-CH=CH-	H	H	-OCH ₂ COOH
942	7-F-Q	-CH=CH-	H	H	-OCH(CH ₃)COOH
943	7-F-Q	-CH=CH-	H	H	-OCH ₂ CH ₂ COOH
944	7-F-Q	-CH=CH-	H	H	-OCH ₂ CH(CH ₃)COOH
945	7-F-Q	-CH=CH-	H	H	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
946	7-F-Q	-CH=CH-	H	H	-SCH ₂ COOH
947	7-F-Q	-CH=CH-	H	H	-SCH(CH ₃)COOH
948	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
949	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)COOH
950	7-F-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₃) ₂ COOH
951	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₂ CH ₃)COOH
952	7-F-Q	-CH=CH-	H	H	-SCH(CH ₃)CH ₂ COOH
953	7-F-Q	-CH=CH-	H	H	-SC(CH ₃) ₂ CH ₂ COOH
954	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CH ₂ COOH
955	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)CH ₂ COOH
956	7-F-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₃) ₂ CH ₂ COOH
957	7-F-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
958	7-F-Q	-CH=CH-	H	H	-SCH ₂ CF ₂ COOH
959	7-F-Q	-CH=CH-	H	H	-SCH ₂ CF ₂ CH ₂ COOH
960	7-F-Q	-CH=CH-	H	H	-SCH ₂ -Tet
961	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ -Tet
962	7-F-Q	-CH=CH-	H	H	-SCH ₂ NHSO ₂ CF ₃
963	7-F-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ CH ₃
964	7-F-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ CF ₃
965	7-F-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ Ph
966	7-F-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ (2-CH ₃ -Ph)
967	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ NHSO ₂ CF ₃
968	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CH ₃
969	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CF ₃
970	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ Ph

971	7-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ (2-CH ₃ -Ph)
972	7-F-Q	-CH=CH-	H	H	-CH ₂ COOH
973	7-F-Q	-CH=CH-	H	H	-CH ₂ CH ₂ COOH
974	7-F-Q	-CH=CH-	H	H	-CH ₂ CH ₂ CH ₂ COOH
975	7-F-Q	-CH=CH-	H	9-F	-OCH ₂ COOH
976	7-F-Q	-CH=CH-	H	9-F	-OCH(CH ₃)COOH
977	7-F-Q	-CH=CH-	H	9-F	-OCH ₂ CH ₂ COOH
978	7-F-Q	-CH=CH-	H	9-F	-SCH ₂ COOH
979	7-F-Q	-CH=CH-	H	9-F	-SCH(CH ₃)COOH
980	7-F-Q	-CH=CH-	H	9-F	-SCH ₂ CH ₂ COOH
981	7-F-Q	-CH=CH-	H	9-F	-SCH ₂ CH(CH ₃)COOH
982	7-F-Q	-CH=CH-	H	9-F	-SCH ₂ CH(CH ₂ CH ₃)COOH
983	7-F-Q	-CH=CH-	H	9-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
984	7-F-Q	-CH=CH-	H	8-F	-OCH ₂ COOH
985	7-F-Q	-CH=CH-	H	8-F	-OCH ₂ CH ₂ COOH
986	7-F-Q	-CH=CH-	H	8-F	-SCH ₂ COOH
987	7-F-Q	-CH=CH-	H	8-F	-SCH ₂ CH ₂ COOH
988	7-F-Q	-CH=CH-	H	8-F	-SCH ₂ CH(CH ₃)COOH
989	7-F-Q	-CH=CH-	H	8-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
990	7-F-Q	-CH=CH-	H	7-F	-OCH ₂ COOH
991	7-F-Q	-CH=CH-	H	7-F	-OCH ₂ CH ₂ COOH
992	7-F-Q	-CH=CH-	H	7-F	-SCH ₂ COOH
993	7-F-Q	-CH=CH-	H	7-F	-SCH ₂ CH ₂ COOH
994	7-F-Q	-CH=CH-	H	7-F	-SCH ₂ CH(CH ₃)COOH
995	7-F-Q	-CH=CH-	H	6-F	-OCH ₂ COOH
996	7-F-Q	-CH=CH-	H	6-F	-OCH(CH ₃)COOH
997	7-F-Q	-CH=CH-	H	6-F	-OCH ₂ CH ₂ COOH
998	7-F-Q	-CH=CH-	H	6-F	-OCH ₂ CH(CH ₃)COOH
999	7-F-Q	-CH=CH-	H	6-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1000	7-F-Q	-CH=CH-	H	6-F	-SCH ₂ COOH
1001	7-F-Q	-CH=CH-	H	6-F	-SCH(CH ₃)COOH
1002	7-F-Q	-CH=CH-	H	6-F	-SCH ₂ CH ₂ COOH
1003	7-F-Q	-CH=CH-	H	6-F	-SCH ₂ CH(CH ₃)COOH

1004	7-F-Q	-CH=CH-	H	6-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1005	7-F-Q	-CH=CH-	H	9-Cl	-OCH ₂ COOH
1006	7-F-Q	-CH=CH-	H	9-Cl	-OCH(CH ₃)COOH
1007	7-F-Q	-CH=CH-	H	9-Cl	-OCH ₂ CH ₂ COOH
1008	7-F-Q	-CH=CH-	H	9-Cl	-OCH ₂ CH(CH ₃)COOH
1009	7-F-Q	-CH=CH-	H	9-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1010	7-F-Q	-CH=CH-	H	9-Cl	-SCH ₂ COOH
1011	7-F-Q	-CH=CH-	H	9-Cl	-SCH(CH ₃)COOH
1012	7-F-Q	-CH=CH-	H	9-Cl	-SCH ₂ CH ₂ COOH
1013	7-F-Q	-CH=CH-	H	9-Cl	-SCH ₂ CH(CH ₃)COOH
1014	7-F-Q	-CH=CH-	H	9-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1015	7-F-Q	-CH=CH-	H	8-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1016	7-F-Q	-CH=CH-	H	8-Cl	-OCH ₂ COOH
1017	7-F-Q	-CH=CH-	H	8-Cl	-OCH ₂ CH ₂ COOH
1018	7-F-Q	-CH=CH-	H	8-Cl	-SCH ₂ COOH
1019	7-F-Q	-CH=CH-	H	8-Cl	-SCH ₂ CH ₂ COOH
1020	7-F-Q	-CH=CH-	H	8-Cl	-SCH ₂ CH(CH ₃)COOH
1021	7-F-Q	-CH=CH-	H	7-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1022	7-F-Q	-CH=CH-	H	7-Cl	-OCH ₂ COOH
1023	7-F-Q	-CH=CH-	H	7-Cl	-OCH ₂ CH ₂ COOH
1024	7-F-Q	-CH=CH-	H	7-Cl	-SCH ₂ COOH
1025	7-F-Q	-CH=CH-	H	7-Cl	-SCH ₂ CH ₂ COOH
1026	7-F-Q	-CH=CH-	H	7-Cl	-SCH ₂ CH(CH ₃)COOH
1027	7-F-Q	-CH=CH-	H	6-Cl	-OCH ₂ COOH
1028	7-F-Q	-CH=CH-	H	6-Cl	-OCH(CH ₃)COOH
1029	7-F-Q	-CH=CH-	H	6-Cl	-OCH ₂ CH ₂ COOH
1030	7-F-Q	-CH=CH-	H	6-Cl	-OCH ₂ CH(CH ₃)COOH
1031	7-F-Q	-CH=CH-	H	6-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1032	7-F-Q	-CH=CH-	H	6-Cl	-SCH ₂ COOH
1033	7-F-Q	-CH=CH-	H	6-Cl	-SCH(CH ₃)COOH
1034	7-F-Q	-CH=CH-	H	6-Cl	-SCH ₂ CH ₂ COOH
1035	7-F-Q	-CH=CH-	H	6-Cl	-SCH ₂ CH(CH ₃)COOH
1036	7-F-Q	-CH=CH-	H	6-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH

1037	7-F-Q	-CH=CH-	H	9-CN	-OCH ₂ COOH
1038	7-F-Q	-CH=CH-	H	9-CN	-OCH(CH ₃)COOH
1039	7-F-Q	-CH=CH-	H	9-CN	-OCH ₂ CH ₂ COOH
1040	7-F-Q	-CH=CH-	H	9-CN	-OCH ₂ CH(CH ₃)COOH
1041	7-F-Q	-CH=CH-	H	9-CN	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1042	7-F-Q	-CH=CH-	H	9-CN	-SCH ₂ COOH
1043	7-F-Q	-CH=CH-	H	9-CN	-SCH(CH ₃)COOH
1044	7-F-Q	-CH=CH-	H	9-CN	-SCH ₂ CH ₂ COOH
1045	7-F-Q	-CH=CH-	H	9-CN	-SCH ₂ CH(CH ₃)COOH
1046	7-F-Q	-CH=CH-	H	9-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1047	7-F-Q	-CH=CH-	H	8-CN	-OCH ₂ COOH
1048	7-F-Q	-CH=CH-	H	8-CN	-OCH ₂ CH ₂ COOH
1049	7-F-Q	-CH=CH-	H	8-CN	-SCH ₂ COOH
1050	7-F-Q	-CH=CH-	H	8-CN	-SCH ₂ CH ₂ COOH
1051	7-F-Q	-CH=CH-	H	8-CN	-SCH ₂ CH(CH ₃)COOH
1052	7-F-Q	-CH=CH-	H	8-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1053	7-F-Q	-CH=CH-	H	7-CN	-OCH ₂ COOH
1054	7-F-Q	-CH=CH-	H	7-CN	-OCH ₂ CH ₂ COOH
1055	7-F-Q	-CH=CH-	H	7-CN	-SCH ₂ COOH
1056	7-F-Q	-CH=CH-	H	7-CN	-SCH ₂ CH ₂ COOH
1057	7-F-Q	-CH=CH-	H	7-CN	-SCH ₂ CH(CH ₃)COOH
1058	7-F-Q	-CH=CH-	H	9-CH ₃	-OCH ₂ COOH
1059	7-F-Q	-CH=CH-	H	9-CH ₃	-OCH ₂ CH ₂ COOH
1060	7-F-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ COOH
1061	7-F-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH ₂ COOH
1062	7-F-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH(CH ₃)COOH
1063	7-F-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH(CH ₂ CH ₃)COOH
1064	7-F-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1065	7-F-Q	-CH=CH-	H	8-CH ₃	-OCH ₂ COOH
1066	7-F-Q	-CH=CH-	H	8-CH ₃	-OCH ₂ CH ₂ COOH
1067	7-F-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ COOH
1068	7-F-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ CH ₂ COOH
1069	7-F-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ CH(CH ₃)COOH

1070	7-F-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1071	7-F-Q	-CH=CH-	H	7-CH ₃	-OCH ₂ COOH
1072	7-F-Q	-CH=CH-	H	7-CH ₃	-OCH ₂ CH ₂ COOH
1073	7-F-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ COOH
1074	7-F-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ CH ₂ COOH
1075	7-F-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ CH(CH ₃)COOH
1076	7-F-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ COOH
1077	7-F-Q	-CH=CH-	H	9-CF ₃	-OCH(CH ₃)COOH
1078	7-F-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ CH ₂ COOH
1079	7-F-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ CH(CH ₃)COOH
1080	7-F-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1081	7-F-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ COOH
1082	7-F-Q	-CH=CH-	H	9-CF ₃	-SCH(CH ₃)COOH
1083	7-F-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ CH ₂ COOH
1084	7-F-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ CH(CH ₃)COOH
1085	7-F-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1086	7-F-Q	-CH=CH-	H	8-CF ₃	-OCH ₂ COOH
1087	7-F-Q	-CH=CH-	H	8-CF ₃	-OCH ₂ CH ₂ COOH
1088	7-F-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ COOH
1089	7-F-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ CH ₂ COOH
1090	7-F-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ CH(CH ₃)COOH
1091	7-F-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1092	7-F-Q	-CH=CH-	H	7-CF ₃	-OCH ₂ COOH
1093	7-F-Q	-CH=CH-	H	7-CF ₃	-OCH ₂ CH ₂ COOH
1094	7-F-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ COOH
1095	7-F-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ CH ₂ COOH
1096	7-F-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ CH(CH ₃)COOH
1097	7-F-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ COOH
1098	7-F-Q	-CH=CH-	H	9-C≡CH	-OCH(CH ₃)COOH
1099	7-F-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ CH ₂ COOH
1100	7-F-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ CH(CH ₃)COOH
1101	7-F-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1102	7-F-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ COOH

1103	7-F-Q	-CH=CH-	H	9-C≡CH	-SCH(CH ₃)COOH
1104	7-F-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ CH ₂ COOH
1105	7-F-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ CH(CH ₃)COOH
1106	7-F-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1107	7-F-Q	-CH=CH-	H	8-C≡CH	-OCH ₂ COOH
1108	7-F-Q	-CH=CH-	H	8-C≡CH	-OCH ₂ CH ₂ COOH
1109	7-F-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ COOH
1110	7-F-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ CH ₂ COOH
1111	7-F-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ CH(CH ₃)COOH
1112	7-F-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1113	7-F-Q	-CH=CH-	H	7-C≡CH	-OCH ₂ COOH
1114	7-F-Q	-CH=CH-	H	7-C≡CH	-OCH ₂ CH ₂ COOH
1115	7-F-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ COOH
1116	7-F-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ CH ₂ COOH
1117	7-F-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ CH(CH ₃)COOH
1118	7-F-Q	-CH=CH-	H	9-CH ₂ OH	-OCH ₂ COOH
1119	7-F-Q	-CH=CH-	H	9-CH ₂ OH	-SCH ₂ CH ₂ COOH
1120	7-F-Q	-CH=CH-	H	8-CH ₂ OH	-OCH ₂ COOH
1121	7-F-Q	-CH=CH-	H	8-CH ₂ OH	-SCH ₂ CH ₂ COOH
1122	7-F-Q	-CH=CH-	H	7-CH ₂ OH	-OCH ₂ COOH
1123	7-F-Q	-CH=CH-	H	7-CH ₂ OH	-SCH ₂ CH ₂ COOH
1124	7-F-Q	-CH=CH-	H	9- C(CH ₃) ₂ OH	-OCH ₂ COOH
1125	7-F-Q	-CH=CH-	H	9- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1126	7-F-Q	-CH=CH-	H	8- C(CH ₃) ₂ OH	-OCH ₂ COOH
1127	7-F-Q	-CH=CH-	H	8- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1128	7-F-Q	-CH=CH-	H	7- C(CH ₃) ₂ OH	-OCH ₂ COOH
1129	7-F-Q	-CH=CH-	H	7- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH

1130	7-F-Q	-CH=CH-	H	9-OCH ₃	-OCH ₂ COOH
1131	7-F-Q	-CH=CH-	H	9-OCH ₃	-SCH ₂ CH ₂ COOH
1132	7-F-Q	-CH=CH-	H	8-OCH ₃	-OCH ₂ COOH
1133	7-F-Q	-CH=CH-	H	8-OCH ₃	-SCH ₂ CH ₂ COOH
1134	7-F-Q	-CH=CH-	H	7-OCH ₃	-OCH ₂ COOH
1135	7-F-Q	-CH=CH-	H	7-OCH ₃	-SCH ₂ CH ₂ COOH
1136	7-F-Q	-CH=CH-	H	9-OCHF ₂	-OCH ₂ COOH
1137	7-F-Q	-CH=CH-	H	9-OCHF ₂	-SCH ₂ CH ₂ COOH
1138	7-F-Q	-CH=CH-	H	8-OCHF ₂	-OCH ₂ COOH
1139	7-F-Q	-CH=CH-	H	8-OCHF ₂	-SCH ₂ CH ₂ COOH
1140	7-F-Q	-CH=CH-	H	7-OCHF ₂	-OCH ₂ COOH
1141	7-F-Q	-CH=CH-	H	7-OCHF ₂	-SCH ₂ CH ₂ COOH
1142	7-F-Q	-CH=CH-	H	9-OCF ₃	-OCH ₂ COOH
1143	7-F-Q	-CH=CH-	H	9-OCF ₃	-SCH ₂ CH ₂ COOH
1144	7-F-Q	-CH=CH-	H	8-OCF ₃	-OCH ₂ COOH
1145	7-F-Q	-CH=CH-	H	8-OCF ₃	-SCH ₂ CH ₂ COOH
1146	7-F-Q	-CH=CH-	H	7-OCF ₃	-OCH ₂ COOH
1147	7-F-Q	-CH=CH-	H	7-OCF ₃	-SCH ₂ CH ₂ COOH
1148	7-F-Q	-CH=CH-	H	9-SOCH ₃	-OCH ₂ COOH
1149	7-F-Q	-CH=CH-	H	9-SOCH ₃	-SCH ₂ CH ₂ COOH
1150	7-F-Q	-CH=CH-	H	8-SOCH ₃	-OCH ₂ COOH
1151	7-F-Q	-CH=CH-	H	8-SOCH ₃	-SCH ₂ CH ₂ COOH
1152	7-F-Q	-CH=CH-	H	7-SOCH ₃	-OCH ₂ COOH
1153	7-F-Q	-CH=CH-	H	7-SOCH ₃	-SCH ₂ CH ₂ COOH
1154	7-F-Q	-CH=CH-	H	9-SO ₂ CH ₃	-OCH ₂ COOH
1155	7-F-Q	-CH=CH-	H	9-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1156	7-F-Q	-CH=CH-	H	8-SO ₂ CH ₃	-OCH ₂ COOH
1157	7-F-Q	-CH=CH-	H	8-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1158	7-F-Q	-CH=CH-	H	7-SO ₂ CH ₃	-OCH ₂ COOH
1159	7-F-Q	-CH=CH-	H	7-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1160	7-F-Q	-CH=CH-	H	9-CH=CH ₂	-SCH ₂ CH ₂ COOH
1161	7-F-Q	-CH=CH-	H	8-CH=CH ₂	-SCH ₂ CH ₂ COOH
1162	7-F-Q	-CH=CH-	H	7-CH=CH ₂	-SCH ₂ CH ₂ COOH

1163	7-F-Q	-CH=CH-	H	9-NO ₂	-OCH ₂ COOH
1164	7-F-Q	-CH=CH-	H	9-NO ₂	-SCH ₂ CH ₂ COOH
1165	7-F-Q	-CH=CH-	H	8-NO ₂	-OCH ₂ COOH
1166	7-F-Q	-CH=CH-	H	8-NO ₂	-SCH ₂ CH ₂ COOH
1167	7-F-Q	-CH=CH-	H	7-NO ₂	-OCH ₂ COOH
1168	7-F-Q	-CH=CH-	H	7-NO ₂	-SCH ₂ CH ₂ COOH
1169	7-F-Q	-CH=CH-	1-F	H	-SCH ₂ CH ₂ COOH
1170	7-F-Q	-CH=CH-	2-F	H	-SCH ₂ CH ₂ COOH
1171	7-F-Q	-CH=CH-	1-Cl	H	-SCH ₂ CH ₂ COOH
1172	7-F-Q	-CH=CH-	2-Cl	H	-SCH ₂ CH ₂ COOH
1173	7-F-Q	-CH=CH-	1-CH ₃	H	-SCH ₂ CH ₂ COOH
1174	7-F-Q	-CH=CH-	2-CH ₃	H	-SCH ₂ CH ₂ COOH
1175	7-F-Q	-CH=CH-	1-OCH ₃	H	-SCH ₂ CH ₂ COOH
1176	7-F-Q	-CH=CH-	2-OCH ₃	H	-SCH ₂ CH ₂ COOH
1177	7-F-Q	-CH=CH-	1-NO ₂	H	-SCH ₂ CH ₂ COOH
1178	7-F-Q	-CH=CH-	2-NO ₂	H	-SCH ₂ CH ₂ COOH
1179	7-F-Q	-CH=CH-	1-CN	H	-SCH ₂ CH ₂ COOH
1180	7-F-Q	-CH=CH-	2-CN	H	-SCH ₂ CH ₂ COOH
1181	7-F-Q	-CH ₂ O-	H	H	-OCH ₂ COOH
1182	7-F-Q	-CH ₂ O-	H	H	-OCH(CH ₃)COOH
1183	7-F-Q	-CH ₂ O-	H	H	-OCH ₂ CH ₂ COOH
1184	7-F-Q	-CH ₂ O-	H	H	-OCH ₂ CH(CH ₃)COOH
1185	7-F-Q	-CH ₂ O-	H	H	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1186	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ COOH
1187	7-F-Q	-CH ₂ O-	H	H	-SCH(CH ₃)COOH
1188	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
1189	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)COOH
1190	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₃) ₂ COOH
1191	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₂ CH ₃)COOH
1192	7-F-Q	-CH ₂ O-	H	H	-SCH(CH ₃)CH ₂ COOH
1193	7-F-Q	-CH ₂ O-	H	H	-SC(CH ₃) ₂ CH ₂ COOH
1194	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CH ₂ COOH
1195	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)CH ₂ COOH

1196	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₃) ₂ CH ₂ COOH
1197	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1198	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CF ₂ COOH
1199	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CF ₂ CH ₂ COOH
1200	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ -Tet
1201	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ -Tet
1202	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ NHSO ₂ CF ₃
1203	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ CH ₃
1204	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ CF ₃
1205	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ Ph
1206	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ (2-CH ₃ -Ph)
1207	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ NHSO ₂ CF ₃
1208	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CH ₃
1209	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CF ₃
1210	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ Ph
1211	7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ (2-CH ₃ -Ph)
1212	7-F-Q	-CH ₂ O-	H	H	-CH ₂ COOH
1213	7-F-Q	-CH ₂ O-	H	H	-CH ₂ CH ₂ COOH
1214	7-F-Q	-CH ₂ O-	H	H	-CH ₂ CH ₂ CH ₂ COOH
1215	7-F-Q	-CH ₂ O-	H	9-F	-OCH ₂ COOH
1216	7-F-Q	-CH ₂ O-	H	9-F	-OCH(CH ₃)COOH
1217	7-F-Q	-CH ₂ O-	H	9-F	-OCH ₂ CH ₂ COOH
1218	7-F-Q	-CH ₂ O-	H	9-F	-OCH ₂ CH(CH ₃)COOH
1219	7-F-Q	-CH ₂ O-	H	9-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1220	7-F-Q	-CH ₂ O-	H	9-F	-SCH ₂ COOH
1221	7-F-Q	-CH ₂ O-	H	9-F	-SCH(CH ₃)COOH
1222	7-F-Q	-CH ₂ O-	H	9-F	-SCH ₂ CH ₂ COOH
1223	7-F-Q	-CH ₂ O-	H	9-F	-SCH ₂ CH(CH ₃)COOH
1224	7-F-Q	-CH ₂ O-	H	9-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1225	7-F-Q	-CH ₂ O-	H	8-F	-OCH ₂ COOH
1226	7-F-Q	-CH ₂ O-	H	8-F	-OCH ₂ CH ₂ COOH
1227	7-F-Q	-CH ₂ O-	H	8-F	-SCH ₂ COOH
1228	7-F-Q	-CH ₂ O-	H	8-F	-SCH ₂ CH ₂ COOH

1229	7-F-Q	-CH ₂ O-	H	8-F	-SCH ₂ CH(CH ₃)COOH
1230	7-F-Q	-CH ₂ O-	H	8-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1231	7-F-Q	-CH ₂ O-	H	7-F	-OCH ₂ COOH
1232	7-F-Q	-CH ₂ O-	H	7-F	-OCH ₂ CH ₂ COOH
1233	7-F-Q	-CH ₂ O-	H	7-F	-SCH ₂ COOH
1234	7-F-Q	-CH ₂ O-	H	7-F	-SCH ₂ CH ₂ COOH
1235	7-F-Q	-CH ₂ O-	H	7-F	-SCH ₂ CH(CH ₃)COOH
1236	7-F-Q	-CH ₂ O-	H	6-F	-OCH ₂ COOH
1237	7-F-Q	-CH ₂ O-	H	6-F	-OCH(CH ₃)COOH
1238	7-F-Q	-CH ₂ O-	H	6-F	-OCH ₂ CH ₂ COOH
1239	7-F-Q	-CH ₂ O-	H	6-F	-OCH ₂ CH(CH ₃)COOH
1240	7-F-Q	-CH ₂ O-	H	6-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1241	7-F-Q	-CH ₂ O-	H	6-F	-SCH ₂ COOH
1242	7-F-Q	-CH ₂ O-	H	6-F	-SCH(CH ₃)COOH
1243	7-F-Q	-CH ₂ O-	H	6-F	-SCH ₂ CH ₂ COOH
1244	7-F-Q	-CH ₂ O-	H	6-F	-SCH ₂ CH(CH ₃)COOH
1245	7-F-Q	-CH ₂ O-	H	6-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1246	7-F-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ COOH
1247	7-F-Q	-CH ₂ O-	H	9-Cl	-OCH(CH ₃)COOH
1248	7-F-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ CH ₂ COOH
1249	7-F-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ CH(CH ₃)COOH
1250	7-F-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1251	7-F-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ COOH
1252	7-F-Q	-CH ₂ O-	H	9-Cl	-SCH(CH ₃)COOH
1253	7-F-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ CH ₂ COOH
1254	7-F-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ CH(CH ₃)COOH
1255	7-F-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1256	7-F-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1257	7-F-Q	-CH ₂ O-	H	8-Cl	-OCH ₂ COOH
1258	7-F-Q	-CH ₂ O-	H	8-Cl	-OCH ₂ CH ₂ COOH
1259	7-F-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ COOH
1260	7-F-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ CH ₂ COOH
1261	7-F-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ CH(CH ₃)COOH

1295	7-F-Q	-CH ₂ O-	H	8-CH ₃	-OCH ₂ COOH
1296	7-F-Q	-CH ₂ O-	H	8-CH ₃	-OCH ₂ CH ₂ COOH
1297	7-F-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ COOH
1298	7-F-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ CH ₂ COOH
1299	7-F-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ CH(CH ₃)COOH
1300	7-F-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1301	7-F-Q	-CH ₂ O-	H	7-CH ₃	-OCH ₂ COOH
1302	7-F-Q	-CH ₂ O-	H	7-CH ₃	-OCH ₂ CH ₂ COOH
1303	7-F-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ COOH
1304	7-F-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ CH ₂ COOH
1305	7-F-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ CH(CH ₃)COOH
1306	7-F-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ COOH
1307	7-F-Q	-CH ₂ O-	H	9-CF ₃	-OCH(CH ₃)COOH
1308	7-F-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ CH ₂ COOH
1309	7-F-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ CH(CH ₃)COOH
1310	7-F-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1311	7-F-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ COOH
1312	7-F-Q	-CH ₂ O-	H	9-CF ₃	-SCH(CH ₃)COOH
1313	7-F-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ CH ₂ COOH
1314	7-F-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ CH(CH ₃)COOH
1315	7-F-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1316	7-F-Q	-CH ₂ O-	H	8-CF ₃	-OCH ₂ COOH
1317	7-F-Q	-CH ₂ O-	H	8-CF ₃	-OCH ₂ CH ₂ COOH
1318	7-F-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ COOH
1319	7-F-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ CH ₂ COOH
1320	7-F-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ CH(CH ₃)COOH
1321	7-F-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1322	7-F-Q	-CH ₂ O-	H	7-CF ₃	-OCH ₂ COOH
1323	7-F-Q	-CH ₂ O-	H	7-CF ₃	-OCH ₂ CH ₂ COOH
1324	7-F-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ COOH
1325	7-F-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ CH ₂ COOH
1326	7-F-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ CH(CH ₃)COOH
1327	7-F-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ COOH

1328	7-F-Q	-CH ₂ O-	H	9-C≡CH	-OCH(CH ₃)COOH
1329	7-F-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ CH ₂ COOH
1330	7-F-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ CH(CH ₃)COOH
1331	7-F-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1332	7-F-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ COOH
1333	7-F-Q	-CH ₂ O-	H	9-C≡CH	-SCH(CH ₃)COOH
1334	7-F-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ CH ₂ COOH
1335	7-F-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ CH(CH ₃)COOH
1336	7-F-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1337	7-F-Q	-CH ₂ O-	H	8-C≡CH	-OCH ₂ COOH
1338	7-F-Q	-CH ₂ O-	H	8-C≡CH	-OCH ₂ CH ₂ COOH
1339	7-F-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ COOH
1340	7-F-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ CH ₂ COOH
1341	7-F-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ CH(CH ₃)COOH
1342	7-F-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1343	7-F-Q	-CH ₂ O-	H	7-C≡CH	-OCH ₂ COOH
1344	7-F-Q	-CH ₂ O-	H	7-C≡CH	-OCH ₂ CH ₂ COOH
1345	7-F-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ COOH
1346	7-F-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ CH ₂ COOH
1347	7-F-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ CH(CH ₃)COOH
1348	7-F-Q	-CH ₂ O-	H	9-CH ₂ OH	-OCH ₂ COOH
1349	7-F-Q	-CH ₂ O-	H	9-CH ₂ OH	-SCH ₂ CH ₂ COOH
1350	7-F-Q	-CH ₂ O-	H	8-CH ₂ OH	-OCH ₂ COOH
1351	7-F-Q	-CH ₂ O-	H	8-CH ₂ OH	-SCH ₂ CH ₂ COOH
1352	7-F-Q	-CH ₂ O-	H	7-CH ₂ OH	-OCH ₂ COOH
1353	7-F-Q	-CH ₂ O-	H	7-CH ₂ OH	-SCH ₂ CH ₂ COOH
1354	7-F-Q	-CH ₂ O-	H	9- C(CH ₃) ₂ OH	-OCH ₂ COOH
1355	7-F-Q	-CH ₂ O-	H	9- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1356	7-F-Q	-CH ₂ O-	H	8- C(CH ₃) ₂ OH	-OCH ₂ COOH

1357	7-F-Q	-CH ₂ O-	H	8- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1358	7-F-Q	-CH ₂ O-	H	7- C(CH ₃) ₂ OH	-OCH ₂ COOH
1359	7-F-Q	-CH ₂ O-	H	7- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1360	7-F-Q	-CH ₂ O-	H	9-OCH ₃	-OCH ₂ COOH
1361	7-F-Q	-CH ₂ O-	H	9-OCH ₃	-SCH ₂ CH ₂ COOH
1362	7-F-Q	-CH ₂ O-	H	8-OCH ₃	-OCH ₂ COOH
1363	7-F-Q	-CH ₂ O-	H	8-OCH ₃	-SCH ₂ CH ₂ COOH
1364	7-F-Q	-CH ₂ O-	H	7-OCH ₃	-OCH ₂ COOH
1365	7-F-Q	-CH ₂ O-	H	7-OCH ₃	-SCH ₂ CH ₂ COOH
1366	7-F-Q	-CH ₂ O-	H	9-OCHF ₂	-OCH ₂ COOH
1367	7-F-Q	-CH ₂ O-	H	9-OCHF ₂	-SCH ₂ CH ₂ COOH
1368	7-F-Q	-CH ₂ O-	H	8-OCHF ₂	-OCH ₂ COOH
1369	7-F-Q	-CH ₂ O-	H	8-OCHF ₂	-SCH ₂ CH ₂ COOH
1370	7-F-Q	-CH ₂ O-	H	7-OCHF ₂	-OCH ₂ COOH
1371	7-F-Q	-CH ₂ O-	H	7-OCHF ₂	-SCH ₂ CH ₂ COOH
1372	7-F-Q	-CH ₂ O-	H	9-OCF ₃	-OCH ₂ COOH
1373	7-F-Q	-CH ₂ O-	H	9-OCF ₃	-SCH ₂ CH ₂ COOH
1374	7-F-Q	-CH ₂ O-	H	8-OCF ₃	-OCH ₂ COOH
1375	7-F-Q	-CH ₂ O-	H	8-OCF ₃	-SCH ₂ CH ₂ COOH
1376	7-F-Q	-CH ₂ O-	H	7-OCF ₃	-OCH ₂ COOH
1377	7-F-Q	-CH ₂ O-	H	7-OCF ₃	-SCH ₂ CH ₂ COOH
1378	7-F-Q	-CH ₂ O-	H	9-SOCH ₃	-OCH ₂ COOH
1379	7-F-Q	-CH ₂ O-	H	9-SOCH ₃	-SCH ₂ CH ₂ COOH
1380	7-F-Q	-CH ₂ O-	H	8-SOCH ₃	-OCH ₂ COOH
1381	7-F-Q	-CH ₂ O-	H	8-SOCH ₃	-SCH ₂ CH ₂ COOH
1382	7-F-Q	-CH ₂ O-	H	7-SOCH ₃	-OCH ₂ COOH
1383	7-F-Q	-CH ₂ O-	H	7-SOCH ₃	-SCH ₂ CH ₂ COOH
1384	7-F-Q	-CH ₂ O-	H	9-SO ₂ CH ₃	-OCH ₂ COOH
1385	7-F-Q	-CH ₂ O-	H	9-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1386	7-F-Q	-CH ₂ O-	H	8-SO ₂ CH ₃	-OCH ₂ COOH

1387	7-F-Q	-CH ₂ O-	H	8-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1388	7-F-Q	-CH ₂ O-	H	7-SO ₂ CH ₃	-OCH ₂ COOH
1389	7-F-Q	-CH ₂ O-	H	7-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1390	7-F-Q	-CH ₂ O-	H	9-CH=CH ₂	-SCH ₂ CH ₂ COOH
1391	7-F-Q	-CH ₂ O-	H	8-CH=CH ₂	-SCH ₂ CH ₂ COOH
1392	7-F-Q	-CH ₂ O-	H	7-CH=CH ₂	-SCH ₂ CH ₂ COOH
1393	7-F-Q	-CH ₂ O-	H	9-NO ₂	-OCH ₂ COOH
1394	7-F-Q	-CH ₂ O-	H	9-NO ₂	-SCH ₂ CH ₂ COOH
1395	7-F-Q	-CH ₂ O-	H	8-NO ₂	-OCH ₂ COOH
1396	7-F-Q	-CH ₂ O-	H	8-NO ₂	-SCH ₂ CH ₂ COOH
1397	7-F-Q	-CH ₂ O-	H	7-NO ₂	-OCH ₂ COOH
1398	7-F-Q	-CH ₂ O-	H	7-NO ₂	-SCH ₂ CH ₂ COOH
1399	7-F-Q	-CH ₂ O-	1-F	H	-SCH ₂ CH ₂ COOH
1400	7-F-Q	-CH ₂ O-	2-F	H	-SCH ₂ CH ₂ COOH
1401	7-F-Q	-CH ₂ O-	1-Cl	H	-SCH ₂ CH ₂ COOH
1402	7-F-Q	-CH ₂ O-	2-Cl	H	-SCH ₂ CH ₂ COOH
1403	7-F-Q	-CH ₂ O-	1-CH ₃	H	-SCH ₂ CH ₂ COOH
1404	7-F-Q	-CH ₂ O-	2-CH ₃	H	-SCH ₂ CH ₂ COOH
1405	7-F-Q	-CH ₂ O-	1-OCH ₃	H	-SCH ₂ CH ₂ COOH
1406	7-F-Q	-CH ₂ O-	2-OCH ₃	H	-SCH ₂ CH ₂ COOH
1407	7-F-Q	-CH ₂ O-	1-NO ₂	H	-SCH ₂ CH ₂ COOH
1408	7-F-Q	-CH ₂ O-	2-NO ₂	H	-SCH ₂ CH ₂ COOH
1409	7-F-Q	-CH ₂ O-	1-CN	H	-SCH ₂ CH ₂ COOH
1410	7-F-Q	-CH ₂ O-	2-CN	H	-SCH ₂ CH ₂ COOH
1411	7-Cl-Q	-CH=CH-	H	H	-OCH ₂ COOH
1412	7-Cl-Q	-CH=CH-	H	H	-OCH(CH ₃)COOH
1413	7-Cl-Q	-CH=CH-	H	H	-OCH ₂ CH ₂ COOH
1414	7-Cl-Q	-CH=CH-	H	H	-OCH ₂ CH(CH ₃)COOH
1415	7-Cl-Q	-CH=CH-	H	H	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1416	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ COOH
1417	7-Cl-Q	-CH=CH-	H	H	-SCH(CH ₃)COOH
1418	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
1419	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)COOH

1420	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₃) ₂ COOH
1421	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₂ CH ₃)COOH
1422	7-Cl-Q	-CH=CH-	H	H	-SCH(CH ₃)CH ₂ COOH
1423	7-Cl-Q	-CH=CH-	H	H	-SC(CH ₃) ₂ CH ₂ COOH
1424	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CH ₂ COOH
1425	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)CH ₂ COOH
1426	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₃) ₂ CH ₂ COOH
1427	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1428	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CF ₂ COOH
1429	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CF ₂ CH ₂ COOH
1430	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ -Tet
1431	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ -Tet
1432	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ NHSO ₂ CF ₃
1433	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ CH ₃
1434	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ CF ₃
1435	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ Ph
1436	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ (2-CH ₃ -Ph)
1437	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ NHSO ₂ CF ₃
1438	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CH ₃
1439	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CF ₃
1440	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ Ph
1441	7-Cl-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ (2-CH ₃ -Ph)
1442	7-Cl-Q	-CH=CH-	H	H	-CH ₂ COOH
1443	7-Cl-Q	-CH=CH-	H	H	-CH ₂ CH ₂ COOH
1444	7-Cl-Q	-CH=CH-	H	H	-CH ₂ CH ₂ CH ₂ COOH
1445	7-Cl-Q	-CH=CH-	H	9-F	-OCH ₂ COOH
1446	7-Cl-Q	-CH=CH-	H	9-F	-OCH(CH ₃)COOH
1447	7-Cl-Q	-CH=CH-	H	9-F	-OCH ₂ CH ₂ COOH
1448	7-Cl-Q	-CH=CH-	H	9-F	-SCH ₂ COOH
1449	7-Cl-Q	-CH=CH-	H	9-F	-SCH(CH ₃)COOH
1450	7-Cl-Q	-CH=CH-	H	9-F	-SCH ₂ CH ₂ COOH
1451	7-Cl-Q	-CH=CH-	H	9-F	-SCH ₂ CH(CH ₃)COOH
1452	7-Cl-Q	-CH=CH-	H	9-F	-SCH ₂ CH(CH ₂ CH ₃)COOH

1453	7-Cl-Q	-CH=CH-	H	9-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1454	7-Cl-Q	-CH=CH-	H	8-F	-OCH ₂ COOH
1455	7-Cl-Q	-CH=CH-	H	8-F	-OCH ₂ CH ₂ COOH
1456	7-Cl-Q	-CH=CH-	H	8-F	-SCH ₂ COOH
1457	7-Cl-Q	-CH=CH-	H	8-F	-SCH ₂ CH ₂ COOH
1458	7-Cl-Q	-CH=CH-	H	8-F	-SCH ₂ CH(CH ₃)COOH
1459	7-Cl-Q	-CH=CH-	H	8-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1460	7-Cl-Q	-CH=CH-	H	7-F	-OCH ₂ COOH
1461	7-Cl-Q	-CH=CH-	H	7-F	-OCH ₂ CH ₂ COOH
1462	7-Cl-Q	-CH=CH-	H	7-F	-SCH ₂ COOH
1463	7-Cl-Q	-CH=CH-	H	7-F	-SCH ₂ CH ₂ COOH
1464	7-Cl-Q	-CH=CH-	H	7-F	-SCH ₂ CH(CH ₃)COOH
1465	7-Cl-Q	-CH=CH-	H	6-F	-OCH ₂ COOH
1466	7-Cl-Q	-CH=CH-	H	6-F	-OCH(CH ₃)COOH
1467	7-Cl-Q	-CH=CH-	H	6-F	-OCH ₂ CH ₂ COOH
1468	7-Cl-Q	-CH=CH-	H	6-F	-OCH ₂ CH(CH ₃)COOH
1469	7-Cl-Q	-CH=CH-	H	6-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1470	7-Cl-Q	-CH=CH-	H	6-F	-SCH ₂ COOH
1471	7-Cl-Q	-CH=CH-	H	6-F	-SCH(CH ₃)COOH
1472	7-Cl-Q	-CH=CH-	H	6-F	-SCH ₂ CH ₂ COOH
1473	7-Cl-Q	-CH=CH-	H	6-F	-SCH ₂ CH(CH ₃)COOH
1474	7-Cl-Q	-CH=CH-	H	6-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1475	7-Cl-Q	-CH=CH-	H	9-Cl	-OCH ₂ COOH
1476	7-Cl-Q	-CH=CH-	H	9-Cl	-OCH(CH ₃)COOH
1477	7-Cl-Q	-CH=CH-	H	9-Cl	-OCH ₂ CH ₂ COOH
1478	7-Cl-Q	-CH=CH-	H	9-Cl	-OCH ₂ CH(CH ₃)COOH
1479	7-Cl-Q	-CH=CH-	H	9-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1480	7-Cl-Q	-CH=CH-	H	9-Cl	-SCH ₂ COOH
1481	7-Cl-Q	-CH=CH-	H	9-Cl	-SCH(CH ₃)COOH
1482	7-Cl-Q	-CH=CH-	H	9-Cl	-SCH ₂ CH ₂ COOH
1483	7-Cl-Q	-CH=CH-	H	9-Cl	-SCH ₂ CH(CH ₃)COOH
1484	7-Cl-Q	-CH=CH-	H	9-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1485	7-Cl-Q	-CH=CH-	H	8-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH

1486	7-Cl-Q	-CH=CH-	H	8-Cl	-OCH ₂ COOH
1487	7-Cl-Q	-CH=CH-	H	8-Cl	-OCH ₂ CH ₂ COOH
1488	7-Cl-Q	-CH=CH-	H	8-Cl	-SCH ₂ COOH
1489	7-Cl-Q	-CH=CH-	H	8-Cl	-SCH ₂ CH ₂ COOH
1490	7-Cl-Q	-CH=CH-	H	8-Cl	-SCH ₂ CH(CH ₃)COOH
1491	7-Cl-Q	-CH=CH-	H	7-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1492	7-Cl-Q	-CH=CH-	H	7-Cl	-OCH ₂ COOH
1493	7-Cl-Q	-CH=CH-	H	7-Cl	-OCH ₂ CH ₂ COOH
1494	7-Cl-Q	-CH=CH-	H	7-Cl	-SCH ₂ COOH
1495	7-Cl-Q	-CH=CH-	H	7-Cl	-SCH ₂ CH ₂ COOH
1496	7-Cl-Q	-CH=CH-	H	7-Cl	-SCH ₂ CH(CH ₃)COOH
1497	7-Cl-Q	-CH=CH-	H	6-Cl	-OCH ₂ COOH
1498	7-Cl-Q	-CH=CH-	H	6-Cl	-OCH(CH ₃)COOH
1499	7-Cl-Q	-CH=CH-	H	6-Cl	-OCH ₂ CH ₂ COOH
1500	7-Cl-Q	-CH=CH-	H	6-Cl	-OCH ₂ CH(CH ₃)COOH
1501	7-Cl-Q	-CH=CH-	H	6-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1502	7-Cl-Q	-CH=CH-	H	6-Cl	-SCH ₂ COOH
1503	7-Cl-Q	-CH=CH-	H	6-Cl	-SCH(CH ₃)COOH
1504	7-Cl-Q	-CH=CH-	H	6-Cl	-SCH ₂ CH ₂ COOH
1505	7-Cl-Q	-CH=CH-	H	6-Cl	-SCH ₂ CH(CH ₃)COOH
1506	7-Cl-Q	-CH=CH-	H	6-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1507	7-Cl-Q	-CH=CH-	H	9-CN	-OCH ₂ COOH
1508	7-Cl-Q	-CH=CH-	H	9-CN	-OCH(CH ₃)COOH
1509	7-Cl-Q	-CH=CH-	H	9-CN	-OCH ₂ CH ₂ COOH
1510	7-Cl-Q	-CH=CH-	H	9-CN	-OCH ₂ CH(CH ₃)COOH
1511	7-Cl-Q	-CH=CH-	H	9-CN	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1512	7-Cl-Q	-CH=CH-	H	9-CN	-SCH ₂ COOH
1513	7-Cl-Q	-CH=CH-	H	9-CN	-SCH(CH ₃)COOH
1514	7-Cl-Q	-CH=CH-	H	9-CN	-SCH ₂ CH ₂ COOH
1515	7-Cl-Q	-CH=CH-	H	9-CN	-SCH ₂ CH(CH ₃)COOH
1516	7-Cl-Q	-CH=CH-	H	9-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1517	7-Cl-Q	-CH=CH-	H	8-CN	-OCH ₂ COOH
1518	7-Cl-Q	-CH=CH-	H	8-CN	-OCH ₂ CH ₂ COOH

1519	7-Cl-Q	-CH=CH-	H	8-CN	-SCH ₂ COOH
1520	7-Cl-Q	-CH=CH-	H	8-CN	-SCH ₂ CH ₂ COOH
1521	7-Cl-Q	-CH=CH-	H	8-CN	-SCH ₂ CH(CH ₃)COOH
1522	7-Cl-Q	-CH=CH-	H	8-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1523	7-Cl-Q	-CH=CH-	H	7-CN	-OCH ₂ COOH
1524	7-Cl-Q	-CH=CH-	H	7-CN	-OCH ₂ CH ₂ COOH
1525	7-Cl-Q	-CH=CH-	H	7-CN	-SCH ₂ COOH
1526	7-Cl-Q	-CH=CH-	H	7-CN	-SCH ₂ CH ₂ COOH
1527	7-Cl-Q	-CH=CH-	H	7-CN	-SCH ₂ CH(CH ₃)COOH
1528	7-Cl-Q	-CH=CH-	H	9-CH ₃	-OCH ₂ COOH
1529	7-Cl-Q	-CH=CH-	H	9-CH ₃	-OCH ₂ CH ₂ COOH
1530	7-Cl-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ COOH
1531	7-Cl-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH ₂ COOH
1532	7-Cl-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH(CH ₃)COOH
1533	7-Cl-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ CH(CH ₂ CH ₃)COOH
1534	7-Cl-Q	-CH=CH-	H	9-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1535	7-Cl-Q	-CH=CH-	H	8-CH ₃	-OCH ₂ COOH
1536	7-Cl-Q	-CH=CH-	H	8-CH ₃	-OCH ₂ CH ₂ COOH
1537	7-Cl-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ COOH
1538	7-Cl-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ CH ₂ COOH
1539	7-Cl-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ CH(CH ₃)COOH
1540	7-Cl-Q	-CH=CH-	H	8-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1541	7-Cl-Q	-CH=CH-	H	7-CH ₃	-OCH ₂ COOH
1542	7-Cl-Q	-CH=CH-	H	7-CH ₃	-OCH ₂ CH ₂ COOH
1543	7-Cl-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ COOH
1544	7-Cl-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ CH ₂ COOH
1545	7-Cl-Q	-CH=CH-	H	7-CH ₃	-SCH ₂ CH(CH ₃)COOH
1546	7-Cl-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ COOH
1547	7-Cl-Q	-CH=CH-	H	9-CF ₃	-OCH(CH ₃)COOH
1548	7-Cl-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ CH ₂ COOH
1549	7-Cl-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ CH(CH ₃)COOH
1550	7-Cl-Q	-CH=CH-	H	9-CF ₃	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1551	7-Cl-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ COOH

1552	7-Cl-Q	-CH=CH-	H	9-CF ₃	-SCH(CH ₃)COOH
1553	7-Cl-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ CH ₂ COOH
1554	7-Cl-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ CH(CH ₃)COOH
1555	7-Cl-Q	-CH=CH-	H	9-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1556	7-Cl-Q	-CH=CH-	H	8-CF ₃	-OCH ₂ COOH
1557	7-Cl-Q	-CH=CH-	H	8-CF ₃	-OCH ₂ CH ₂ COOH
1558	7-Cl-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ COOH
1559	7-Cl-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ CH ₂ COOH
1560	7-Cl-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ CH(CH ₃)COOH
1561	7-Cl-Q	-CH=CH-	H	8-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1562	7-Cl-Q	-CH=CH-	H	7-CF ₃	-OCH ₂ COOH
1563	7-Cl-Q	-CH=CH-	H	7-CF ₃	-OCH ₂ CH ₂ COOH
1564	7-Cl-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ COOH
1565	7-Cl-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ CH ₂ COOH
1566	7-Cl-Q	-CH=CH-	H	7-CF ₃	-SCH ₂ CH(CH ₃)COOH
1567	7-Cl-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ COOH
1568	7-Cl-Q	-CH=CH-	H	9-C≡CH	-OCH(CH ₃)COOH
1569	7-Cl-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ CH ₂ COOH
1570	7-Cl-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ CH(CH ₃)COOH
1571	7-Cl-Q	-CH=CH-	H	9-C≡CH	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1572	7-Cl-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ COOH
1573	7-Cl-Q	-CH=CH-	H	9-C≡CH	-SCH(CH ₃)COOH
1574	7-Cl-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ CH ₂ COOH
1575	7-Cl-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ CH(CH ₃)COOH
1576	7-Cl-Q	-CH=CH-	H	9-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1577	7-Cl-Q	-CH=CH-	H	8-C≡CH	-OCH ₂ COOH
1578	7-Cl-Q	-CH=CH-	H	8-C≡CH	-OCH ₂ CH ₂ COOH
1579	7-Cl-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ COOH
1580	7-Cl-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ CH ₂ COOH
1581	7-Cl-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ CH(CH ₃)COOH
1582	7-Cl-Q	-CH=CH-	H	8-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1583	7-Cl-Q	-CH=CH-	H	7-C≡CH	-OCH ₂ COOH
1584	7-Cl-Q	-CH=CH-	H	7-C≡CH	-OCH ₂ CH ₂ COOH

1585	7-Cl-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ COOH
1586	7-Cl-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ CH ₂ COOH
1587	7-Cl-Q	-CH=CH-	H	7-C≡CH	-SCH ₂ CH(CH ₃)COOH
1588	7-Cl-Q	-CH=CH-	H	9-CH ₂ OH	-OCH ₂ COOH
1589	7-Cl-Q	-CH=CH-	H	9-CH ₂ OH	-SCH ₂ CH ₂ COOH
1590	7-Cl-Q	-CH=CH-	H	8-CH ₂ OH	-OCH ₂ COOH
1591	7-Cl-Q	-CH=CH-	H	8-CH ₂ OH	-SCH ₂ CH ₂ COOH
1592	7-Cl-Q	-CH=CH-	H	7-CH ₂ OH	-OCH ₂ COOH
1593	7-Cl-Q	-CH=CH-	H	7-CH ₂ OH	-SCH ₂ CH ₂ COOH
1594	7-Cl-Q	-CH=CH-	H	9- C(CH ₃) ₂ OH	-OCH ₂ COOH
1595	7-Cl-Q	-CH=CH-	H	9- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1596	7-Cl-Q	-CH=CH-	H	8- C(CH ₃) ₂ OH	-OCH ₂ COOH
1597	7-Cl-Q	-CH=CH-	H	8- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1598	7-Cl-Q	-CH=CH-	H	7- C(CH ₃) ₂ OH	-OCH ₂ COOH
1599	7-Cl-Q	-CH=CH-	H	7- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1600	7-Cl-Q	-CH=CH-	H	9-OCH ₃	-OCH ₂ COOH
1601	7-Cl-Q	-CH=CH-	H	9-OCH ₃	-SCH ₂ CH ₂ COOH
1602	7-Cl-Q	-CH=CH-	H	8-OCH ₃	-OCH ₂ COOH
1603	7-Cl-Q	-CH=CH-	H	8-OCH ₃	-SCH ₂ CH ₂ COOH
1604	7-Cl-Q	-CH=CH-	H	7-OCH ₃	-OCH ₂ COOH
1605	7-Cl-Q	-CH=CH-	H	7-OCH ₃	-SCH ₂ CH ₂ COOH
1606	7-Cl-Q	-CH=CH-	H	9-OCHF ₂	-OCH ₂ COOH
1607	7-Cl-Q	-CH=CH-	H	9-OCHF ₂	-SCH ₂ CH ₂ COOH
1608	7-Cl-Q	-CH=CH-	H	8-OCHF ₂	-OCH ₂ COOH
1609	7-Cl-Q	-CH=CH-	H	8-OCHF ₂	-SCH ₂ CH ₂ COOH
1610	7-Cl-Q	-CH=CH-	H	7-OCHF ₂	-OCH ₂ COOH
1611	7-Cl-Q	-CH=CH-	H	7-OCHF ₂	-SCH ₂ CH ₂ COOH

1612	7-Cl-Q	-CH=CH-	H	9-OCF ₃	-OCH ₂ COOH
1613	7-Cl-Q	-CH=CH-	H	9-OCF ₃	-SCH ₂ CH ₂ COOH
1614	7-Cl-Q	-CH=CH-	H	8-OCF ₃	-OCH ₂ COOH
1615	7-Cl-Q	-CH=CH-	H	8-OCF ₃	-SCH ₂ CH ₂ COOH
1616	7-Cl-Q	-CH=CH-	H	7-OCF ₃	-OCH ₂ COOH
1617	7-Cl-Q	-CH=CH-	H	7-OCF ₃	-SCH ₂ CH ₂ COOH
1618	7-Cl-Q	-CH=CH-	H	9-SOCH ₃	-OCH ₂ COOH
1619	7-Cl-Q	-CH=CH-	H	9-SOCH ₃	-SCH ₂ CH ₂ COOH
1620	7-Cl-Q	-CH=CH-	H	8-SOCH ₃	-OCH ₂ COOH
1621	7-Cl-Q	-CH=CH-	H	8-SOCH ₃	-SCH ₂ CH ₂ COOH
1622	7-Cl-Q	-CH=CH-	H	7-SOCH ₃	-OCH ₂ COOH
1623	7-Cl-Q	-CH=CH-	H	7-SOCH ₃	-SCH ₂ CH ₂ COOH
1624	7-Cl-Q	-CH=CH-	H	9-SO ₂ CH ₃	-OCH ₂ COOH
1625	7-Cl-Q	-CH=CH-	H	9-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1626	7-Cl-Q	-CH=CH-	H	8-SO ₂ CH ₃	-OCH ₂ COOH
1627	7-Cl-Q	-CH=CH-	H	8-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1628	7-Cl-Q	-CH=CH-	H	7-SO ₂ CH ₃	-OCH ₂ COOH
1629	7-Cl-Q	-CH=CH-	H	7-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1630	7-Cl-Q	-CH=CH-	H	9-CH=CH ₂	-SCH ₂ CH ₂ COOH
1631	7-Cl-Q	-CH=CH-	H	8-CH=CH ₂	-SCH ₂ CH ₂ COOH
1632	7-Cl-Q	-CH=CH-	H	7-CH=CH ₂	-SCH ₂ CH ₂ COOH
1633	7-Cl-Q	-CH=CH-	H	9-NO ₂	-OCH ₂ COOH
1634	7-Cl-Q	-CH=CH-	H	9-NO ₂	-SCH ₂ CH ₂ COOH
1635	7-Cl-Q	-CH=CH-	H	8-NO ₂	-OCH ₂ COOH
1636	7-Cl-Q	-CH=CH-	H	8-NO ₂	-SCH ₂ CH ₂ COOH
1637	7-Cl-Q	-CH=CH-	H	7-NO ₂	-OCH ₂ COOH
1638	7-Cl-Q	-CH=CH-	H	7-NO ₂	-SCH ₂ CH ₂ COOH
1639	7-Cl-Q	-CH=CH-	1-F	H	-SCH ₂ CH ₂ COOH
1640	7-Cl-Q	-CH=CH-	2-F	H	-SCH ₂ CH ₂ COOH
1641	7-Cl-Q	-CH=CH-	1-Cl	H	-SCH ₂ CH ₂ COOH
1642	7-Cl-Q	-CH=CH-	2-Cl	H	-SCH ₂ CH ₂ COOH
1643	7-Cl-Q	-CH=CH-	1-CH ₃	H	-SCH ₂ CH ₂ COOH
1644	7-Cl-Q	-CH=CH-	2-CH ₃	H	-SCH ₂ CH ₂ COOH

1645	7-Cl-Q	-CH=CH-	1-OCH ₃	H	-SCH ₂ CH ₂ COOH
1646	7-Cl-Q	-CH=CH-	2-OCH ₃	H	-SCH ₂ CH ₂ COOH
1647	7-Cl-Q	-CH=CH-	1-NO ₂	H	-SCH ₂ CH ₂ COOH
1648	7-Cl-Q	-CH=CH-	2-NO ₂	H	-SCH ₂ CH ₂ COOH
1649	7-Cl-Q	-CH=CH-	1-CN	H	-SCH ₂ CH ₂ COOH
1650	7-Cl-Q	-CH=CH-	2-CN	H	-SCH ₂ CH ₂ COOH
1651	7-Cl-Q	-CH ₂ O-	H	H	-OCH ₂ COOH
1652	7-Cl-Q	-CH ₂ O-	H	H	-OCH(CH ₃)COOH
1653	7-Cl-Q	-CH ₂ O-	H	H	-OCH ₂ CH ₂ COOH
1654	7-Cl-Q	-CH ₂ O-	H	H	-OCH ₂ CH(CH ₃)COOH
1655	7-Cl-Q	-CH ₂ O-	H	H	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1656	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ COOH
1657	7-Cl-Q	-CH ₂ O-	H	H	-SCH(CH ₃)COOH
1658	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
1659	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)COOH
1660	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₃) ₂ COOH
1661	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₂ CH ₃)COOH
1662	7-Cl-Q	-CH ₂ O-	H	H	-SCH(CH ₃)CH ₂ COOH
1663	7-Cl-Q	-CH ₂ O-	H	H	-SC(CH ₃) ₂ CH ₂ COOH
1664	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CH ₂ COOH
1665	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)CH ₂ COOH
1666	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₃) ₂ CH ₂ COOH
1667	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1668	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CF ₂ COOH
1669	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CF ₂ CH ₂ COOH
1670	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ -Tet
1671	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ -Tet
1672	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ NHSO ₂ CF ₃
1673	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ CH ₃
1674	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ CF ₃
1675	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ Ph
1676	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ (2-CH ₃ -Ph)
1677	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ NHSO ₂ CF ₃

1678	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CH ₃
1679	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CF ₃
1680	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ Ph
1681	7-Cl-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ (2-CH ₃ -Ph)
1682	7-Cl-Q	-CH ₂ O-	H	H	-CH ₂ COOH
1683	7-Cl-Q	-CH ₂ O-	H	H	-CH ₂ CH ₂ COOH
1684	7-Cl-Q	-CH ₂ O-	H	H	-CH ₂ CH ₂ CH ₂ COOH
1685	7-Cl-Q	-CH ₂ O-	H	9-F	-OCH ₂ COOH
1686	7-Cl-Q	-CH ₂ O-	H	9-F	-OCH(CH ₃)COOH
1687	7-Cl-Q	-CH ₂ O-	H	9-F	-OCH ₂ CH ₂ COOH
1688	7-Cl-Q	-CH ₂ O-	H	9-F	-OCH ₂ CH(CH ₃)COOH
1689	7-Cl-Q	-CH ₂ O-	H	9-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1690	7-Cl-Q	-CH ₂ O-	H	9-F	-SCH ₂ COOH
1691	7-Cl-Q	-CH ₂ O-	H	9-F	-SCH(CH ₃)COOH
1692	7-Cl-Q	-CH ₂ O-	H	9-F	-SCH ₂ CH ₂ COOH
1693	7-Cl-Q	-CH ₂ O-	H	9-F	-SCH ₂ CH(CH ₃)COOH
1694	7-Cl-Q	-CH ₂ O-	H	9-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1695	7-Cl-Q	-CH ₂ O-	H	8-F	-OCH ₂ COOH
1696	7-Cl-Q	-CH ₂ O-	H	8-F	-OCH ₂ CH ₂ COOH
1697	7-Cl-Q	-CH ₂ O-	H	8-F	-SCH ₂ COOH
1698	7-Cl-Q	-CH ₂ O-	H	8-F	-SCH ₂ CH ₂ COOH
1699	7-Cl-Q	-CH ₂ O-	H	8-F	-SCH ₂ CH(CH ₃)COOH
1700	7-Cl-Q	-CH ₂ O-	H	8-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1701	7-Cl-Q	-CH ₂ O-	H	7-F	-OCH ₂ COOH
1702	7-Cl-Q	-CH ₂ O-	H	7-F	-OCH ₂ CH ₂ COOH
1703	7-Cl-Q	-CH ₂ O-	H	7-F	-SCH ₂ COOH
1704	7-Cl-Q	-CH ₂ O-	H	7-F	-SCH ₂ CH ₂ COOH
1705	7-Cl-Q	-CH ₂ O-	H	7-F	-SCH ₂ CH(CH ₃)COOH
1706	7-Cl-Q	-CH ₂ O-	H	6-F	-OCH ₂ COOH
1707	7-Cl-Q	-CH ₂ O-	H	6-F	-OCH(CH ₃)COOH
1708	7-Cl-Q	-CH ₂ O-	H	6-F	-OCH ₂ CH ₂ COOH
1709	7-Cl-Q	-CH ₂ O-	H	6-F	-OCH ₂ CH(CH ₃)COOH
1710	7-Cl-Q	-CH ₂ O-	H	6-F	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH

1711	7-Cl-Q	-CH ₂ O-	H	6-F	-SCH ₂ COOH
1712	7-Cl-Q	-CH ₂ O-	H	6-F	-SCH(CH ₃)COOH
1713	7-Cl-Q	-CH ₂ O-	H	6-F	-SCH ₂ CH ₂ COOH
1714	7-Cl-Q	-CH ₂ O-	H	6-F	-SCH ₂ CH(CH ₃)COOH
1715	7-Cl-Q	-CH ₂ O-	H	6-F	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1716	7-Cl-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ COOH
1717	7-Cl-Q	-CH ₂ O-	H	9-Cl	-OCH(CH ₃)COOH
1718	7-Cl-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ CH ₂ COOH
1719	7-Cl-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ CH(CH ₃)COOH
1720	7-Cl-Q	-CH ₂ O-	H	9-Cl	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1721	7-Cl-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ COOH
1722	7-Cl-Q	-CH ₂ O-	H	9-Cl	-SCH(CH ₃)COOH
1723	7-Cl-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ CH ₂ COOH
1724	7-Cl-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ CH(CH ₃)COOH
1725	7-Cl-Q	-CH ₂ O-	H	9-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1726	7-Cl-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1727	7-Cl-Q	-CH ₂ O-	H	8-Cl	-OCH ₂ COOH
1728	7-Cl-Q	-CH ₂ O-	H	8-Cl	-OCH ₂ CH ₂ COOH
1729	7-Cl-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ COOH
1730	7-Cl-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ CH ₂ COOH
1731	7-Cl-Q	-CH ₂ O-	H	8-Cl	-SCH ₂ CH(CH ₃)COOH
1732	7-Cl-Q	-CH ₂ O-	H	7-Cl	-OCH ₂ COOH
1733	7-Cl-Q	-CH ₂ O-	H	7-Cl	-OCH ₂ CH ₂ COOH
1734	7-Cl-Q	-CH ₂ O-	H	7-Cl	-SCH ₂ COOH
1735	7-Cl-Q	-CH ₂ O-	H	7-Cl	-SCH ₂ CH ₂ COOH
1736	7-Cl-Q	-CH ₂ O-	H	7-Cl	-SCH ₂ CH(CH ₃)COOH
1737	7-Cl-Q	-CH ₂ O-	H	9-CN	-OCH ₂ COOH
1738	7-Cl-Q	-CH ₂ O-	H	9-CN	-OCH(CH ₃)COOH
1739	7-Cl-Q	-CH ₂ O-	H	9-CN	-OCH ₂ CH ₂ COOH
1740	7-Cl-Q	-CH ₂ O-	H	9-CN	-OCH ₂ CH(CH ₃)COOH
1741	7-Cl-Q	-CH ₂ O-	H	9-CN	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1742	7-Cl-Q	-CH ₂ O-	H	9-CN	-SCH ₂ COOH
1743	7-Cl-Q	-CH ₂ O-	H	9-CN	-SCH(CH ₃)COOH

1744	7-Cl-Q	-CH ₂ O-	H	9-CN	-SCH ₂ CH ₂ COOH
1745	7-Cl-Q	-CH ₂ O-	H	9-CN	-SCH ₂ CH(CH ₃)COOH
1746	7-Cl-Q	-CH ₂ O-	H	9-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1747	7-Cl-Q	-CH ₂ O-	H	8-CN	-OCH ₂ COOH
1748	7-Cl-Q	-CH ₂ O-	H	8-CN	-OCH ₂ CH ₂ COOH
1749	7-Cl-Q	-CH ₂ O-	H	8-CN	-SCH ₂ COOH
1750	7-Cl-Q	-CH ₂ O-	H	8-CN	-SCH ₂ CH ₂ COOH
1751	7-Cl-Q	-CH ₂ O-	H	8-CN	-SCH ₂ CH(CH ₃)COOH
1752	7-Cl-Q	-CH ₂ O-	H	8-CN	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1753	7-Cl-Q	-CH ₂ O-	H	7-CN	-OCH ₂ COOH
1754	7-Cl-Q	-CH ₂ O-	H	7-CN	-OCH ₂ CH ₂ COOH
1755	7-Cl-Q	-CH ₂ O-	H	7-CN	-SCH ₂ COOH
1756	7-Cl-Q	-CH ₂ O-	H	7-CN	-SCH ₂ CH ₂ COOH
1757	7-Cl-Q	-CH ₂ O-	H	7-CN	-SCH ₂ CH(CH ₃)COOH
1758	7-Cl-Q	-CH ₂ O-	H	9-CH ₃	-OCH ₂ COOH
1759	7-Cl-Q	-CH ₂ O-	H	9-CH ₃	-OCH ₂ CH ₂ COOH
1760	7-Cl-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ COOH
1761	7-Cl-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ CH ₂ COOH
1762	7-Cl-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ CH(CH ₃)COOH
1763	7-Cl-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ CH(CH ₂ CH ₃)COOH
1764	7-Cl-Q	-CH ₂ O-	H	9-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1765	7-Cl-Q	-CH ₂ O-	H	8-CH ₃	-OCH ₂ COOH
1766	7-Cl-Q	-CH ₂ O-	H	8-CH ₃	-OCH ₂ CH ₂ COOH
1767	7-Cl-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ COOH
1768	7-Cl-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ CH ₂ COOH
1769	7-Cl-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ CH(CH ₃)COOH
1770	7-Cl-Q	-CH ₂ O-	H	8-CH ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1771	7-Cl-Q	-CH ₂ O-	H	7-CH ₃	-OCH ₂ COOH
1772	7-Cl-Q	-CH ₂ O-	H	7-CH ₃	-OCH ₂ CH ₂ COOH
1773	7-Cl-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ COOH
1774	7-Cl-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ CH ₂ COOH
1775	7-Cl-Q	-CH ₂ O-	H	7-CH ₃	-SCH ₂ CH(CH ₃)COOH
1776	7-Cl-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ COOH

1777	7-Cl-Q	-CH ₂ O-	H	9-CF ₃	-OCH(CH ₃)COOH
1778	7-Cl-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ CH ₂ COOH
1779	7-Cl-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ CH(CH ₃)COOH
1780	7-Cl-Q	-CH ₂ O-	H	9-CF ₃	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1781	7-Cl-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ COOH
1782	7-Cl-Q	-CH ₂ O-	H	9-CF ₃	-SCH(CH ₃)COOH
1783	7-Cl-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ CH ₂ COOH
1784	7-Cl-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ CH(CH ₃)COOH
1785	7-Cl-Q	-CH ₂ O-	H	9-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1786	7-Cl-Q	-CH ₂ O-	H	8-CF ₃	-OCH ₂ COOH
1787	7-Cl-Q	-CH ₂ O-	H	8-CF ₃	-OCH ₂ CH ₂ COOH
1788	7-Cl-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ COOH
1789	7-Cl-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ CH ₂ COOH
1790	7-Cl-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ CH(CH ₃)COOH
1791	7-Cl-Q	-CH ₂ O-	H	8-CF ₃	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1792	7-Cl-Q	-CH ₂ O-	H	7-CF ₃	-OCH ₂ COOH
1793	7-Cl-Q	-CH ₂ O-	H	7-CF ₃	-OCH ₂ CH ₂ COOH
1794	7-Cl-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ COOH
1795	7-Cl-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ CH ₂ COOH
1796	7-Cl-Q	-CH ₂ O-	H	7-CF ₃	-SCH ₂ CH(CH ₃)COOH
1797	7-Cl-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ COOH
1798	7-Cl-Q	-CH ₂ O-	H	9-C≡CH	-OCH(CH ₃)COOH
1799	7-Cl-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ CH ₂ COOH
1800	7-Cl-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ CH(CH ₃)COOH
1801	7-Cl-Q	-CH ₂ O-	H	9-C≡CH	-OCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1802	7-Cl-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ COOH
1803	7-Cl-Q	-CH ₂ O-	H	9-C≡CH	-SCH(CH ₃)COOH
1804	7-Cl-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ CH ₂ COOH
1805	7-Cl-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ CH(CH ₃)COOH
1806	7-Cl-Q	-CH ₂ O-	H	9-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1807	7-Cl-Q	-CH ₂ O-	H	8-C≡CH	-OCH ₂ COOH
1808	7-Cl-Q	-CH ₂ O-	H	8-C≡CH	-OCH ₂ CH ₂ COOH
1809	7-Cl-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ COOH

1810	7-Cl-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ CH ₂ COOH
1811	7-Cl-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ CH(CH ₃)COOH
1812	7-Cl-Q	-CH ₂ O-	H	8-C≡CH	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1813	7-Cl-Q	-CH ₂ O-	H	7-C≡CH	-OCH ₂ COOH
1814	7-Cl-Q	-CH ₂ O-	H	7-C≡CH	-OCH ₂ CH ₂ COOH
1815	7-Cl-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ COOH
1816	7-Cl-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ CH ₂ COOH
1817	7-Cl-Q	-CH ₂ O-	H	7-C≡CH	-SCH ₂ CH(CH ₃)COOH
1818	7-Cl-Q	-CH ₂ O-	H	9-CH ₂ OH	-OCH ₂ COOH
1819	7-Cl-Q	-CH ₂ O-	H	9-CH ₂ OH	-SCH ₂ CH ₂ COOH
1820	7-Cl-Q	-CH ₂ O-	H	8-CH ₂ OH	-OCH ₂ COOH
1821	7-Cl-Q	-CH ₂ O-	H	8-CH ₂ OH	-SCH ₂ CH ₂ COOH
1822	7-Cl-Q	-CH ₂ O-	H	7-CH ₂ OH	-OCH ₂ COOH
1823	7-Cl-Q	-CH ₂ O-	H	7-CH ₂ OH	-SCH ₂ CH ₂ COOH
1824	7-Cl-Q	-CH ₂ O-	H	9- C(CH ₃) ₂ OH	-OCH ₂ COOH
1825	7-Cl-Q	-CH ₂ O-	H	9- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1826	7-Cl-Q	-CH ₂ O-	H	8- C(CH ₃) ₂ OH	-OCH ₂ COOH
1827	7-Cl-Q	-CH ₂ O-	H	8- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1828	7-Cl-Q	-CH ₂ O-	H	7- C(CH ₃) ₂ OH	-OCH ₂ COOH
1829	7-Cl-Q	-CH ₂ O-	H	7- C(CH ₃) ₂ OH	-SCH ₂ CH ₂ COOH
1830	7-Cl-Q	-CH ₂ O-	H	9-OCH ₃	-OCH ₂ COOH
1831	7-Cl-Q	-CH ₂ O-	H	9-OCH ₃	-SCH ₂ CH ₂ COOH
1832	7-Cl-Q	-CH ₂ O-	H	8-OCH ₃	-OCH ₂ COOH
1833	7-Cl-Q	-CH ₂ O-	H	8-OCH ₃	-SCH ₂ CH ₂ COOH
1834	7-Cl-Q	-CH ₂ O-	H	7-OCH ₃	-OCH ₂ COOH
1835	7-Cl-Q	-CH ₂ O-	H	7-OCH ₃	-SCH ₂ CH ₂ COOH
1836	7-Cl-Q	-CH ₂ O-	H	9-OCHF ₂	-OCH ₂ COOH

1837	7-Cl-Q	-CH ₂ O-	H	9-OCHF ₂	-SCH ₂ CH ₂ COOH
1838	7-Cl-Q	-CH ₂ O-	H	8-OCHF ₂	-OCH ₂ COOH
1839	7-Cl-Q	-CH ₂ O-	H	8-OCHF ₂	-SCH ₂ CH ₂ COOH
1840	7-Cl-Q	-CH ₂ O-	H	7-OCHF ₂	-OCH ₂ COOH
1841	7-Cl-Q	-CH ₂ O-	H	7-OCHF ₂	-SCH ₂ CH ₂ COOH
1842	7-Cl-Q	-CH ₂ O-	H	9-OCF ₃	-OCH ₂ COOH
1843	7-Cl-Q	-CH ₂ O-	H	9-OCF ₃	-SCH ₂ CH ₂ COOH
1844	7-Cl-Q	-CH ₂ O-	H	8-OCF ₃	-OCH ₂ COOH
1845	7-Cl-Q	-CH ₂ O-	H	8-OCF ₃	-SCH ₂ CH ₂ COOH
1846	7-Cl-Q	-CH ₂ O-	H	7-OCF ₃	-OCH ₂ COOH
1847	7-Cl-Q	-CH ₂ O-	H	7-OCF ₃	-SCH ₂ CH ₂ COOH
1848	7-Cl-Q	-CH ₂ O-	H	9-SOCH ₃	-OCH ₂ COOH
1849	7-Cl-Q	-CH ₂ O-	H	9-SOCH ₃	-SCH ₂ CH ₂ COOH
1850	7-Cl-Q	-CH ₂ O-	H	8-SOCH ₃	-OCH ₂ COOH
1851	7-Cl-Q	-CH ₂ O-	H	8-SOCH ₃	-SCH ₂ CH ₂ COOH
1852	7-Cl-Q	-CH ₂ O-	H	7-SOCH ₃	-OCH ₂ COOH
1853	7-Cl-Q	-CH ₂ O-	H	7-SOCH ₃	-SCH ₂ CH ₂ COOH
1854	7-Cl-Q	-CH ₂ O-	H	9-SO ₂ CH ₃	-OCH ₂ COOH
1855	7-Cl-Q	-CH ₂ O-	H	9-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1856	7-Cl-Q	-CH ₂ O-	H	8-SO ₂ CH ₃	-OCH ₂ COOH
1857	7-Cl-Q	-CH ₂ O-	H	8-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1858	7-Cl-Q	-CH ₂ O-	H	7-SO ₂ CH ₃	-OCH ₂ COOH
1859	7-Cl-Q	-CH ₂ O-	H	7-SO ₂ CH ₃	-SCH ₂ CH ₂ COOH
1860	7-Cl-Q	-CH ₂ O-	H	9-CH=CH ₂	-SCH ₂ CH ₂ COOH
1861	7-Cl-Q	-CH ₂ O-	H	8-CH=CH ₂	-SCH ₂ CH ₂ COOH
1862	7-Cl-Q	-CH ₂ O-	H	7-CH=CH ₂	-SCH ₂ CH ₂ COOH
1863	7-Cl-Q	-CH ₂ O-	H	9-NO ₂	-OCH ₂ COOH
1864	7-Cl-Q	-CH ₂ O-	H	9-NO ₂	-SCH ₂ CH ₂ COOH
1865	7-Cl-Q	-CH ₂ O-	H	8-NO ₂	-OCH ₂ COOH
1866	7-Cl-Q	-CH ₂ O-	H	8-NO ₂	-SCH ₂ CH ₂ COOH
1867	7-Cl-Q	-CH ₂ O-	H	7-NO ₂	-OCH ₂ COOH
1868	7-Cl-Q	-CH ₂ O-	H	7-NO ₂	-SCH ₂ CH ₂ COOH
1869	7-Cl-Q	-CH ₂ O-	1-F	H	-SCH ₂ CH ₂ COOH

1870	7-Cl-Q	-CH ₂ O-	2-F	H	-SCH ₂ CH ₂ COOH
1871	7-Cl-Q	-CH ₂ O-	1-Cl	H	-SCH ₂ CH ₂ COOH
1872	7-Cl-Q	-CH ₂ O-	2-Cl	H	-SCH ₂ CH ₂ COOH
1873	7-Cl-Q	-CH ₂ O-	1-CH ₃	H	-SCH ₂ CH ₂ COOH
1874	7-Cl-Q	-CH ₂ O-	2-CH ₃	H	-SCH ₂ CH ₂ COOH
1875	7-Cl-Q	-CH ₂ O-	1-OCH ₃	H	-SCH ₂ CH ₂ COOH
1876	7-Cl-Q	-CH ₂ O-	2-OCH ₃	H	-SCH ₂ CH ₂ COOH
1877	7-Cl-Q	-CH ₂ O-	1-NO ₂	H	-SCH ₂ CH ₂ COOH
1878	7-Cl-Q	-CH ₂ O-	2-NO ₂	H	-SCH ₂ CH ₂ COOH
1879	7-Cl-Q	-CH ₂ O-	1-CN	H	-SCH ₂ CH ₂ COOH
1880	7-Cl-Q	-CH ₂ O-	2-CN	H	-SCH ₂ CH ₂ COOH
1881	TQ	-CH=CH-	H	H	-OCH ₂ COOH
1882	TQ	-CH=CH-	H	H	-OCH ₂ CH ₂ COOH
1883	TQ	-CH=CH-	H	H	-OCH ₂ CH(CH ₃)COOH
1884	TQ	-CH=CH-	H	H	-SCH ₂ COOH
1885	TQ	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
1886	TQ	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)COOH
1887	TQ	-CH=CH-	H	H	-SCH ₂ C(CH ₃) ₂ COOH
1888	TQ	-CH=CH-	H	H	-SCH ₂ CH(CH ₂ CH ₃)COOH
1889	TQ	-CH=CH-	H	H	-SCH(CH ₃)CH ₂ COOH
1890	TQ	-CH=CH-	H	H	-SC(CH ₃) ₂ CH ₂ COOH
1891	TQ	-CH=CH-	H	H	-SCH ₂ CH ₂ CH ₂ COOH
1892	TQ	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)CH ₂ COOH
1893	TQ	-CH=CH-	H	H	-SCH ₂ C(CH ₃) ₂ CH ₂ COOH
1894	TQ	-CH=CH-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1895	TQ	-CH=CH-	H	H	-SCH ₂ CF ₂ COOH
1896	TQ	-CH=CH-	H	H	-SCH ₂ CF ₂ CH ₂ COOH
1897	TQ	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ CH ₃
1898	TQ	-CH=CH-	H	H	-SCH ₂ CONHSO ₂ CF ₃
1899	TQ	-CH=CH-	H	H	-SCH ₂ CH ₂ NHSO ₂ CF ₃
1900	TQ	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CH ₃
1901	TQ	-CH=CH-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CF ₃
1902	TQ	-CH=CH-	H	H	-CH ₂ CH ₂ COOH

1903	TQ	-CH=CH-	H	H	-CH ₂ CH ₂ CH ₂ COOH
1904	TQ	-CH=CH-	H	9-F	-OCH ₂ COOH
1905	TQ	-CH=CH-	H	9-F	-OCH ₂ CH ₂ COOH
1906	TQ	-CH=CH-	H	9-F	-SCH ₂ COOH
1907	TQ	-CH=CH-	H	9-F	-SCH ₂ CH ₂ COOH
1908	TQ	-CH=CH-	H	8-F	-OCH ₂ COOH
1909	TQ	-CH=CH-	H	8-F	-OCH ₂ CH ₂ COOH
1910	TQ	-CH=CH-	H	8-F	-SCH ₂ COOH
1911	TQ	-CH=CH-	H	8-F	-SCH ₂ CH ₂ COOH
1912	TQ	-CH=CH-	H	9-CN	-OCH ₂ COOH
1913	TQ	-CH=CH-	H	9-CN	-OCH ₂ CH ₂ COOH
1914	TQ	-CH=CH-	H	9-CN	-SCH ₂ COOH
1915	TQ	-CH=CH-	H	9-CN	-SCH ₂ CH ₂ COOH
1916	TQ	-CH=CH-	H	8-CN	-OCH ₂ COOH
1917	TQ	-CH=CH-	H	8-CN	-OCH ₂ CH ₂ COOH
1918	TQ	-CH=CH-	H	8-CN	-SCH ₂ COOH
1919	TQ	-CH=CH-	H	8-CN	-SCH ₂ CH ₂ COOH
1920	TQ	-CH=CH-	H	9-CF ₃	-OCH ₂ COOH
1921	TQ	-CH=CH-	H	9-CF ₃	-SCH ₂ CH ₂ COOH
1922	TQ	-CH=CH-	H	8-CF ₃	-OCH ₂ COOH
1923	TQ	-CH=CH-	H	8-CF ₃	-SCH ₂ CH ₂ COOH
1924	TQ	-CH=CH-	H	9-C≡CH	-OCH ₂ COOH
1925	TQ	-CH=CH-	H	9-C≡CH	-SCH ₂ CH ₂ COOH
1926	TQ	-CH=CH-	H	8-C≡CH	-SCH ₂ CH ₂ COOH
1927	TQ	-CH ₂ O-	H	H	-OCH ₂ COOH
1928	TQ	-CH ₂ O-	H	H	-OCH ₂ CH ₂ COOH
1929	TQ	-CH ₂ O-	H	H	-OCH ₂ CH(CH ₃)COOH
1930	TQ	-CH ₂ O-	H	H	-SCH ₂ COOH
1931	TQ	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
1932	TQ	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)COOH
1933	TQ	-CH ₂ O-	H	H	-SCH ₂ C(CH ₃) ₂ COOH
1934	TQ	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₂ CH ₃)COOH
1935	TQ	-CH ₂ O-	H	H	-SCH(CH ₃)CH ₂ COOH

1936	TQ	-CH ₂ O-	H	H	-SC(CH ₃) ₂ CH ₂ COOH
1937	TQ	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CH ₂ COOH
1938	TQ	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)CH ₂ COOH
1939	TQ	-CH ₂ O-	H	H	-SCH ₂ C(CH ₃) ₂ CH ₂ COOH
1940	TQ	-CH ₂ O-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1941	TQ	-CH ₂ O-	H	H	-SCH ₂ CF ₂ COOH
1942	TQ	-CH ₂ O-	H	H	-SCH ₂ CF ₂ CH ₂ COOH
1943	TQ	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ CH ₃
1944	TQ	-CH ₂ O-	H	H	-SCH ₂ CONHSO ₂ CF ₃
1945	TQ	-CH ₂ O-	H	H	-SCH ₂ CH ₂ NHSO ₂ CF ₃
1946	TQ	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CH ₃
1947	TQ	-CH ₂ O-	H	H	-SCH ₂ CH ₂ CONHSO ₂ CF ₃
1948	TQ	-CH ₂ O-	H	H	-CH ₂ CH ₂ COOH
1949	TQ	-CH ₂ O-	H	H	-CH ₂ CH ₂ CH ₂ COOH
1950	TQ	-CH ₂ O-	H	9-F	-OCH ₂ COOH
1951	TQ	-CH ₂ O-	H	9-F	-OCH ₂ CH ₂ COOH
1952	TQ	-CH ₂ O-	H	9-F	-SCH ₂ COOH
1953	TQ	-CH ₂ O-	H	9-F	-SCH ₂ CH ₂ COOH
1954	TQ	-CH ₂ O-	H	8-F	-OCH ₂ COOH
1955	TQ	-CH ₂ O-	H	8-F	-OCH ₂ CH ₂ COOH
1956	TQ	-CH ₂ O-	H	8-F	-SCH ₂ COOH
1957	TQ	-CH ₂ O-	H	8-F	-SCH ₂ CH ₂ COOH
1958	TQ	-CH ₂ O-	H	9-CN	-OCH ₂ COOH
1959	TQ	-CH ₂ O-	H	9-CN	-OCH ₂ CH ₂ COOH
1960	TQ	-CH ₂ O-	H	9-CN	-SCH ₂ COOH
1961	TQ	-CH ₂ O-	H	9-CN	-SCH ₂ CH ₂ COOH
1962	TQ	-CH ₂ O-	H	8-CN	-OCH ₂ COOH
1963	TQ	-CH ₂ O-	H	8-CN	-OCH ₂ CH ₂ COOH
1964	TQ	-CH ₂ O-	H	8-CN	-SCH ₂ COOH
1965	TQ	-CH ₂ O-	H	8-CN	-SCH ₂ CH ₂ COOH
1966	TQ	-CH ₂ O-	H	9-CF ₃	-OCH ₂ COOH
1967	TQ	-CH ₂ O-	H	9-CF ₃	-SCH ₂ CH ₂ COOH
1968	TQ	-CH ₂ O-	H	8-CF ₃	-OCH ₂ COOH

1969	TQ	-CH ₂ O-	H	8-CF ₃	-SCH ₂ CH ₂ COOH
1970	TQ	-CH ₂ O-	H	9-C≡CH	-OCH ₂ COOH
1971	TQ	-CH ₂ O-	H	9-C≡CH	-SCH ₂ CH ₂ COOH
1972	TQ	-CH ₂ O-	H	8-C≡CH	-SCH ₂ CH ₂ COOH
1973	7-CF ₃ -Q	-CH=CH-	H	H	-OCH ₂ COOH
1974	7-CF ₃ -Q	-CH=CH-	H	H	-OCH ₂ CH ₂ COOH
1975	7-CF ₃ -Q	-CH=CH-	H	H	-SCH ₂ COOH
1976	7-CF ₃ -Q	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
1977	7-CF ₃ -Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)COOH
1978	7-CF ₃ -Q	-CH=CH-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1979	7-CF ₃ -Q	-CH ₂ O-	H	H	-OCH ₂ COOH
1980	7-CF ₃ -Q	-CH ₂ O-	H	H	-OCH ₂ CH ₂ COOH
1981	7-CF ₃ -Q	-CH ₂ O-	H	H	-SCH ₂ COOH
1982	7-CF ₃ -Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
1983	7-CF ₃ -Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)COOH
1984	7-CF ₃ -Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1985	6-F,7-CF ₃ -Q	-CH=CH-	H	H	-OCH ₂ COOH
1986	6-F,7-CF ₃ -Q	-CH=CH-	H	H	-OCH ₂ CH ₂ COOH
1987	6-F,7-CF ₃ -Q	-CH=CH-	H	H	-SCH ₂ COOH
1988	6-F,7-CF ₃ -Q	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
1989	6-F,7-CF ₃ -Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)COOH
1990	6-F,7-CF ₃ -Q	-CH=CH-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1991	6-F,7-CF ₃ -Q	-CH ₂ O-	H	H	-OCH ₂ COOH
1992	6-F,7-CF ₃ -Q	-CH ₂ O-	H	H	-OCH ₂ CH ₂ COOH
1993	6-F,7-CF ₃ -Q	-CH ₂ O-	H	H	-SCH ₂ COOH
1994	6-F,7-CF ₃ -Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
1995	6-F,7-CF ₃ -Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)COOH
1996	6-F,7-CF ₃ -Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
1997	6,7-diCl-Q	-CH=CH-	H	H	-OCH ₂ COOH
1998	6,7-diCl-Q	-CH=CH-	H	H	-OCH ₂ CH ₂ COOH
1999	6,7-diCl-Q	-CH=CH-	H	H	-SCH ₂ COOH
2000	6,7-diCl-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
2001	6,7-diCl-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)COOH

2002	6,7-diCl-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
2003	6,7-diCl-Q	-CH ₂ O-	H	H	-OCH ₂ COOH
2004	6,7-diCl-Q	-CH ₂ O-	H	H	-OCH ₂ CH ₂ COOH
2005	6,7-diCl-Q	-CH ₂ O-	H	H	-SCH ₂ COOH
2006	6,7-diCl-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
2007	6,7-diCl-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)COOH
2008	6,7-diCl-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
2009	6-Cl,7-F-Q	-CH=CH-	H	H	-OCH ₂ COOH
2010	6-Cl,7-F-Q	-CH=CH-	H	H	-OCH ₂ CH ₂ COOH
2011	6-Cl,7-F-Q	-CH=CH-	H	H	-SCH ₂ COOH
2012	6-Cl,7-F-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
2013	6-Cl,7-F-Q	-CH=CH-	H	H	-SCH ₂ CH(CH ₃)COOH
2014	6-Cl,7-F-Q	-CH=CH-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
2015	6-Cl,7-F-Q	-CH ₂ O-	H	H	-OCH ₂ COOH
2016	6-Cl,7-F-Q	-CH ₂ O-	H	H	-OCH ₂ CH ₂ COOH
2017	6-Cl,7-F-Q	-CH ₂ O-	H	H	-SCH ₂ COOH
2018	6-Cl,7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
2019	6-Cl,7-F-Q	-CH ₂ O-	H	H	-SCH ₂ CH(CH ₃)COOH
2020	6-Cl,7-F-Q	-CH ₂ O-	H	H	-SCH ₂ C(CH ₂ CH ₂)CH ₂ COOH
2021	5,6,7-triF-Q	-CH=CH-	H	H	-OCH ₂ COOH
2022	5,6,7-triF-Q	-CH=CH-	H	H	-OCH ₂ CH ₂ COOH
2023	5,6,7-triF-Q	-CH=CH-	H	H	-SCH ₂ COOH
2024	5,6,7-triF-Q	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
2025	5,6,7-triF-Q	-CH ₂ O-	H	H	-OCH ₂ COOH
2026	5,6,7-triF-Q	-CH ₂ O-	H	H	-OCH ₂ CH ₂ COOH
2027	5,6,7-triF-Q	-CH ₂ O-	H	H	-SCH ₂ COOH
2028	5,6,7-triF-Q	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
2029	4-t-Bu-T	-CH=CH-	H	H	-OCH ₂ COOH
2030	4-t-Bu-T	-CH=CH-	H	H	-SCH ₂ COOH
2031	4-t-Bu-T	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
2032	4-t-Bu-T	-CH ₂ O-	H	H	-OCH ₂ COOH
2033	4-t-Bu-T	-CH ₂ O-	H	H	-SCH ₂ COOH
2034	4-t-Bu-T	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH

2035	5-F-BT	-CH=CH-	H	H	-OCH ₂ COOH
2036	5-F-BT	-CH=CH-	H	H	-SCH ₂ COOH
2037	5-F-BT	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
2038	5-F-BT	-CH ₂ O-	H	H	-OCH ₂ COOH
2039	5-F-BT	-CH ₂ O-	H	H	-SCH ₂ COOH
2040	5-F-BT	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
2041	5,6-diF-BT	-CH=CH-	H	H	-OCH ₂ COOH
2042	5,6-diF-BT	-CH=CH-	H	H	-SCH ₂ COOH
2043	5,6-diF-BT	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
2044	5,6-diF-BT	-CH ₂ O-	H	H	-OCH ₂ COOH
2045	5,6-diF-BT	-CH ₂ O-	H	H	-SCH ₂ COOH
2046	5,6-diF-BT	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
2047	6-t-Bu-Py	-CH=CH-	H	H	-OCH ₂ COOH
2048	6-t-Bu-Py	-CH=CH-	H	H	-SCH ₂ COOH
2049	6-t-Bu-Py	-CH=CH-	H	H	-SCH ₂ CH ₂ COOH
2050	6-t-Bu-Py	-CH ₂ O-	H	H	-OCH ₂ COOH
2051	6-t-Bu-Py	-CH ₂ O-	H	H	-SCH ₂ COOH
2052	6-t-Bu-Py	-CH ₂ O-	H	H	-SCH ₂ CH ₂ COOH
2053	6,7-diF-Q	-CH ₂ CH ₂ -	H	H	-OCH ₂ COOH
2054	6,7-diF-Q	-CH ₂ CH ₂ -	H	H	-SCH ₂ COOH
2055	6,7-diF-Q	-CH ₂ CH ₂ -	H	H	-SCH ₂ CH ₂ COOH
2056	6,7-diF-Q	-CH ₂ S-	H	H	-OCH ₂ COOH
2057	6,7-diF-Q	-CH ₂ S-	H	H	-SCH ₂ COOH
2058	6,7-diF-Q	-CH ₂ S-	H	H	-SCH ₂ CH ₂ COOH
2059	6,7-diF-Q	-OCH ₂ -	H	H	-OCH ₂ COOH
2060	6,7-diF-Q	-OCH ₂ -	H	H	-SCH ₂ COOH
2061	6,7-diF-Q	-OCH ₂ -	H	H	-SCH ₂ CH ₂ COOH
2062	6,7-diF-Q	-SCH ₂ -	H	H	-OCH ₂ COOH
2063	6,7-diF-Q	-SCH ₂ -	H	H	-SCH ₂ COOH
2064	6,7-diF-Q	-SCH ₂ -	H	H	-SCH ₂ CH ₂ COOH
2065	7-Cl,6-F-Q	-CH ₂ CH ₂ -	H	H	-OCH ₂ COOH
2066	7-Cl,6-F-Q	-CH ₂ CH ₂ -	H	H	-SCH ₂ COOH
2067	7-Cl,6-F-Q	-CH ₂ CH ₂ -	H	H	-SCH ₂ CH ₂ COOH

2068	7-Cl,6-F-Q	-CH ₂ S-	H	H	-OCH ₂ COOH
2069	7-Cl,6-F-Q	-CH ₂ S-	H	H	-SCH ₂ COOH
2070	7-Cl,6-F-Q	-CH ₂ S-	H	H	-SCH ₂ CH ₂ COOH
2071	7-Cl,6-F-Q	-OCH ₂ -	H	H	-OCH ₂ COOH
2072	7-Cl,6-F-Q	-OCH ₂ -	H	H	-SCH ₂ COOH
2073	7-Cl,6-F-Q	-OCH ₂ -	H	H	-SCH ₂ CH ₂ COOH
2074	7-Cl,6-F-Q	-SCH ₂ -	H	H	-OCH ₂ COOH
2075	7-Cl,6-F-Q	-SCH ₂ -	H	H	-SCH ₂ COOH
2076	7-Cl,6-F-Q	-SCH ₂ -	H	H	-SCH ₂ CH ₂ COOH
2077	7-F-Q	-CH ₂ CH ₂ -	H	H	-OCH ₂ COOH
2078	7-F-Q	-CH ₂ CH ₂ -	H	H	-SCH ₂ COOH
2079	7-F-Q	-CH ₂ CH ₂ -	H	H	-SCH ₂ CH ₂ COOH
2080	7-F-Q	-CH ₂ S-	H	H	-OCH ₂ COOH
2081	7-F-Q	-CH ₂ S-	H	H	-SCH ₂ COOH
2082	7-F-Q	-CH ₂ S-	H	H	-SCH ₂ CH ₂ COOH
2083	7-F-Q	-OCH ₂ -	H	H	-OCH ₂ COOH
2084	7-F-Q	-OCH ₂ -	H	H	-SCH ₂ COOH
2085	7-F-Q	-OCH ₂ -	H	H	-SCH ₂ CH ₂ COOH
2086	7-F-Q	-SCH ₂ -	H	H	-OCH ₂ COOH
2087	7-F-Q	-SCH ₂ -	H	H	-SCH ₂ COOH
2088	7-F-Q	-SCH ₂ -	H	H	-SCH ₂ CH ₂ COOH
2089	7-Cl-Q	-CH ₂ CH ₂ -	H	H	-OCH ₂ COOH
2090	7-Cl-Q	-CH ₂ CH ₂ -	H	H	-SCH ₂ COOH
2091	7-Cl-Q	-CH ₂ CH ₂ -	H	H	-SCH ₂ CH ₂ COOH
2092	7-Cl-Q	-CH ₂ S-	H	H	-OCH ₂ COOH
2093	7-Cl-Q	-CH ₂ S-	H	H	-SCH ₂ COOH
2094	7-Cl-Q	-CH ₂ S-	H	H	-SCH ₂ CH ₂ COOH
2095	7-Cl-Q	-OCH ₂ -	H	H	-OCH ₂ COOH
2096	7-Cl-Q	-OCH ₂ -	H	H	-SCH ₂ COOH
2097	7-Cl-Q	-OCH ₂ -	H	H	-SCH ₂ CH ₂ COOH
2098	7-Cl-Q	-SCH ₂ -	H	H	-OCH ₂ COOH
2099	7-Cl-Q	-SCH ₂ -	H	H	-SCH ₂ COOH
2100	7-Cl-Q	-SCH ₂ -	H	H	-SCH ₂ CH ₂ COOH

Incidentally, in the above-mentioned table, the respective abbreviations mean

t-Bu: t-butyl group, BT: 2-benzothiazolyl group, Tet: 1H-tetrazol-5-yl group, Ph: a phenyl group, Py: 2-pyridyl
 5 group, Q: quinolin-2-yl group, T: 2-thiazolyl group, and TQ: 5,6,7,8-tetrahydroquinolin-2-yl group,

and the $\text{CH}_2\text{C}(\text{CH}_2\text{CH}_2)\text{CH}_2$ portion in the $-\text{OCH}_2\text{C}(\text{CH}_2\text{CH}_2)-\text{CH}_2\text{COOH}$ or $-\text{SCH}_2\text{C}(\text{CH}_2\text{CH}_2)\text{CH}_2\text{COOH}$ represents the formula (a) wherein $o=1$, $p=1$ and $q=1$, and the $\text{CH}_2\text{C}(\text{CH}_2\text{CH}_2)$ portion in
 10 the $-\text{SCH}_2\text{C}(\text{CH}_2\text{CH}_2)\text{COOH}$ represents the formula (a) wherein $o=1$, $p=0$ and $q=1$. Also, in the above-mentioned Table 1, the numeral(s) in the formula (I) represents a substituted position(s) of the substituent(s).

As the more preferred compounds, there may be
 15 mentioned compounds of Compounds No. 1, 2, 3, 4, 5, 6, 7, 8, 9, 11, 12, 13, 14, 15, 16, 17, 24, 25, 27, 33, 34, 35, 36, 37, 38, 39, 40, 41, 42, 43, 44, 47, 49, 55, 56, 57, 58, 59, 60, 61, 62, 63, 64, 65, 66, 67, 68, 69, 70, 71, 72, 73, 74, 76, 78, 79, 87, 89, 92, 94, 97, 99, 102, 103, 104, 105,
 20 106, 107, 110, 112, 118, 121, 122, 123, 124, 128, 136, 138, 139, 140, 141, 142, 143, 144, 145, 146, 149, 157, 159, 160, 161, 162, 163, 164, 165, 166, 167, 170, 179, 184, 185, 190, 191, 196, 197, 202, 203, 220, 223, 224, 241, 243, 244, 245, 246, 247, 248, 249, 250, 251, 257, 264, 273, 274, 275, 277,
 25 280, 281, 282, 283, 284, 285, 288, 296, 298, 299, 300, 301, 302, 303, 304, 305, 306, 308, 309, 310, 311, 312, 313, 314, 315, 319, 320, 327, 329, 331, 332, 333, 334, 335, 336, 337, 340, 348, 351, 366, 368, 371, 372, 373, 374, 375, 376, 378, 379, 387, 392, 393, 394, 396, 397, 400, 420, 421, 426, 427,
 30 429, 432, 433, 435, 453, 454, 455, 456, 471, 472, 473, 474, 475, 476, 477, 478, 479, 481, 482, 483, 484, 485, 486, 487, 494, 495, 497, 503, 504, 505, 506, 507, 508, 509, 510, 511, 512, 513, 514, 517, 519, 525, 526, 527, 528, 529, 530, 531, 532, 533, 534, 535, 536, 537, 538, 539, 540, 541, 542, 543,
 35 544, 546, 548, 549, 557, 558, 559, 562, 564, 567, 569, 572, 573, 574, 575, 576, 577, 580, 582, 588, 591, 592, 593, 594,

598, 606, 608, 609, 610, 611, 612, 613, 614, 615, 616, 619,
627, 629, 630, 631, 632, 633, 634, 635, 636, 637, 640, 649,
654, 655, 660, 661, 666, 667, 672, 673, 690, 693, 694, 711,
713, 714, 715, 716, 717, 718, 719, 720, 721, 727, 734, 743,
5 744, 745, 747, 750, 751, 752, 753, 754, 755, 758, 766, 768,
769, 770, 771, 772, 773, 774, 775, 776, 778, 779, 780, 781,
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805, 806, 807, 810, 818, 821, 836, 838, 841, 842, 843, 844,
845, 846, 848, 849, 857, 862, 863, 864, 866, 867, 870, 890,
10 891, 896, 897, 899, 902, 903, 905, 923, 924, 925, 926, 941,
942, 943, 944, 945, 946, 947, 948, 949, 951, 952, 953, 954,
955, 956, 957, 964, 965, 967, 973, 974, 975, 976, 977, 978,
979, 980, 981, 982, 983, 984, 987, 989, 995, 996, 997, 998,
999, 1000, 1001, 1002, 1003, 1004, 1005, 1006, 1007, 1008,
15 1009, 1010, 1011, 1012, 1013, 1014, 1016, 1018, 1019, 1027,
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1078, 1079, 1080, 1081, 1082, 1083, 1084, 1085, 1086, 1089,
1097, 1099, 1100, 1101, 1102, 1103, 1104, 1105, 1106, 1107,
20 1110, 1119, 1124, 1125, 1130, 1131, 1136, 1137, 1142, 1143,
1160, 1163, 1164, 1181, 1183, 1184, 1185, 1186, 1187, 1188,
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1372, 1373, 1375, 1393, 1394, 1395, 1396, 1411, 1412, 1413,
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1447, 1448, 1449, 1450, 1451, 1452, 1453, 1454, 1457, 1459,
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1475, 1476, 1477, 1478, 1479, 1480, 1481, 1482, 1483, 1484,
35 1486, 1488, 1489, 1497, 1499, 1502, 1504, 1507, 1509, 1512,
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1533, 1534, 1538, 1546, 1548, 1549, 1550, 1551, 1552, 1553,
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1574, 1575, 1576, 1577, 1580, 1589, 1594, 1595, 1600, 1601,
1606, 1607, 1612, 1613, 1630, 1633, 1634, 1651, 1653, 1654,
5 1655, 1656, 1657, 1658, 1659, 1660, 1661, 1667, 1674, 1683,
1684, 1685, 1687, 1690, 1691, 1692, 1693, 1694, 1695, 1698,
1706, 1708, 1709, 1710, 1711, 1712, 1713, 1714, 1715, 1716,
1718, 1719, 1720, 1721, 1722, 1723, 1724, 1725, 1729, 1730,
1737, 1739, 1741, 1742, 1743, 1744, 1745, 1746, 1747, 1750,
10 1758, 1761, 1776, 1778, 1781, 1782, 1783, 1784, 1785, 1786,
1788, 1789, 1797, 1802, 1803, 1804, 1806, 1807, 1810, 1830,
1831, 1836, 1837, 1839, 1842, 1843, 1845, 1863, 1864, 1865,
1866, 1884, 1885, 1886, 1888, 1894, 1903, 1907, 1911, 1915,
1921, 1925, 1931, 1932, 1934, 1940, 1949, 1953, 1961, 1967,
15 1971, 1976, 1978, 1980, 1982, 1984, 1988, 1990, 1994, 1995,
1996, 2000, 2001, 2002, 2004, 2006, 2007, 2008, 2012, 2013,
2014, 2018, 2019, 2020, 2024, 2028, 2031, 2034, 2037, 2040,
2043, 2046, 2049, 2052, 2055, 2058, 2061, 2064, 2067, 2070,
2073, 2076, 2079, 2082, 2085, 2088, 2091, 2094, 2097 or
20 2100,

more preferably compounds are Compounds No. 1, 2, 3,
4, 6, 7, 8, 9, 11, 14, 15, 17, 35, 37, 38, 40, 41, 42, 43,
47, 49, 55, 57, 60, 62, 63, 64, 65, 66, 67, 70, 72, 73, 74,
78, 79, 87, 89, 92, 94, 97, 102, 104, 105, 106, 110, 121,
25 122, 123, 136, 138, 141, 143, 144, 145, 149, 157, 159, 162,
163, 164, 165, 166, 170, 191, 197, 202, 203, 220, 224, 241,
243, 246, 247, 248, 249, 251, 257, 275, 280, 282, 283, 284,
288, 296, 298, 301, 303, 304, 305, 306, 308, 311, 312, 313,
314, 315, 320, 327, 332, 334, 335, 336, 340, 351, 366, 371,
30 372, 373, 374, 375, 376, 379, 387, 393, 394, 400, 420, 421,
426, 427, 432, 433, 453, 456, 471, 473, 474, 476, 477, 478,
479, 481, 484, 485, 487, 505, 507, 508, 510, 511, 512, 513,
517, 519, 525, 527, 530, 532, 533, 534, 535, 536, 537, 540,
542, 543, 544, 548, 549, 557, 558, 559, 562, 564, 567, 572,
35 574, 575, 576, 580, 591, 592, 593, 606, 608, 611, 613, 614,
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672, 673, 690, 694, 711, 713, 716, 717, 718, 719, 721, 727,
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 776, 778, 781, 782, 783, 784, 785, 790, 797, 802, 804, 805,
 806, 810, 821, 836, 841, 842, 843, 844, 845, 846, 849, 857,
 5 863, 864, 870, 890, 891, 896, 897, 902, 903, 923, 926, 941,
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 1425, 1427, 1445, 1447, 1448, 1450, 1451, 1452, 1453, 1457,
 20 1459, 1465, 1467, 1470, 1472, 1473, 1474, 1475, 1476, 1477,
 1480, 1482, 1483, 1484, 1488, 1489, 1497, 1499, 1502, 1504,
 1507, 1512, 1514, 1515, 1516, 1520, 1531, 1532, 1533, 1546,
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 1574, 1575, 1576, 1580, 1601, 1607, 1612, 1613, 1630, 1634,
 25 1651, 1653, 1656, 1657, 1658, 1659, 1661, 1667, 1685, 1690,
 1692, 1693, 1694, 1695, 1698, 1706, 1708, 1711, 1713, 1714,
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 1742, 1744, 1745, 1746, 1750, 1761, 1776, 1781, 1782, 1783,
 1784, 1785, 1786, 1789, 1797, 1803, 1804, 1810, 1830, 1831,
 30 1836, 1837, 1842, 1843, 1863, 1866, 1885, 1886, 1888, 1894,
 1907, 1911, 1915, 1921, 1931, 1932, 1934, 1953, 1961, 1967
 or 1971,

further more preferably compounds of Compounds No.

1, 2, 3, 4, 6, 8, 9, 11, 15, 17, 38, 40, 41, 42, 43, 47,
 35 62, 63, 64, 72, 73, 74, 79, 87, 89, 92, 94, 97, 104, 106,
 110, 121, 136, 143, 144, 145, 149, 157, 163, 164, 165, 166,

191, 197, 203, 220, 247, 248, 249, 251, 257, 282, 283, 284,
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 5 542, 543, 544, 549, 557, 558, 559, 562, 564, 567, 574, 576,
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 10 978, 980, 981, 982, 983, 987, 1002, 1003, 1004, 1012, 1013,
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 1497, 1499, 1502, 1504, 1507, 1514, 1516, 1520, 1531, 1546,
 1553, 1554, 1555, 1559, 1567, 1573, 1574, 1575, 1576, 1601,
 20 1607, 1613, 1630, 1657, 1658, 1659, 1661, 1667, 1692, 1693,
 1694, 1713, 1722, 1723, 1724, 1725, 1744, 1782, 1783, 1784,
 1785, 1803, 1804, 1831, 1837, 1843, 1885, 1886, 1894, 1907
 or 1921.

It is particularly preferably mentioned

- 25 Compound No.1; [3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethen-
 yl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid,
 Compound No.4; 3-{[3-[(E)-2-(6,7-difluoroquinolin-2-yl)-
 ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}-2-methyl-
 propionic acid,
 30 Compound No.8; 3-{[3-[(E)-2-(6,7-difluoroquinolin-2-yl)-
 ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic
 acid,
 Compound No.9; 3-{[3-[(E)-2-(6,7-difluoroquinolin-2-yl)-
 ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}-2-methyl-
 35 propionic acid,
 Compound No.471; [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-

- yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid,
Compound No.473; 3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}propionic
acid,
- 5 Compound No.474; 3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}-2-methyl-
propionic acid,
Compound No.476; [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thioacetic
10 acid,
Compound No.477; 2-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic
acid,
Compound No.478; 3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
15 yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic
acid,
Compound No.487; 3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thiomethyl}-
cyclopropane acetic acid,
- 20 Compound No.505; [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-yl]oxy-
acetic acid,
Compound No.510; 3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-yl]thio}-
25 propionic acid,
Compound No.535; [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-9-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]oxy-
acetic acid,
Compound No.542; 3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
30 yl)ethenyl]-9-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thio}-
propionic acid,
Compound No.549; 3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-8-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thio}-
propionic acid,
- 35 Compound No.557; [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]oxy-

acetic acid,

Compound No.558; 2-{{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}propionic acid,

5 Compound No.559; 3-{{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}-propionic acid,

Compound No.562; [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thio-
10 acetic acid,

Compound No.564; 3-{{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thio}-propionic acid,

Compound No.613; 3-{{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

Compound No.718; 3-{{[3-(7-chloro-6-fluoroquinolin-2-yl)-methoxy-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

20 Compound No.1411; [3-[(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid,

Compound No.1416; [3-[(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thioacetic acid,

Compound No.1418; 3-{{[3-[(E)-2-(7-chloroquinolin-2-yl)-ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

Compound No.1885; 3-{{[3-[(E)-2-(5,6,7,8-tetrahydroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid or

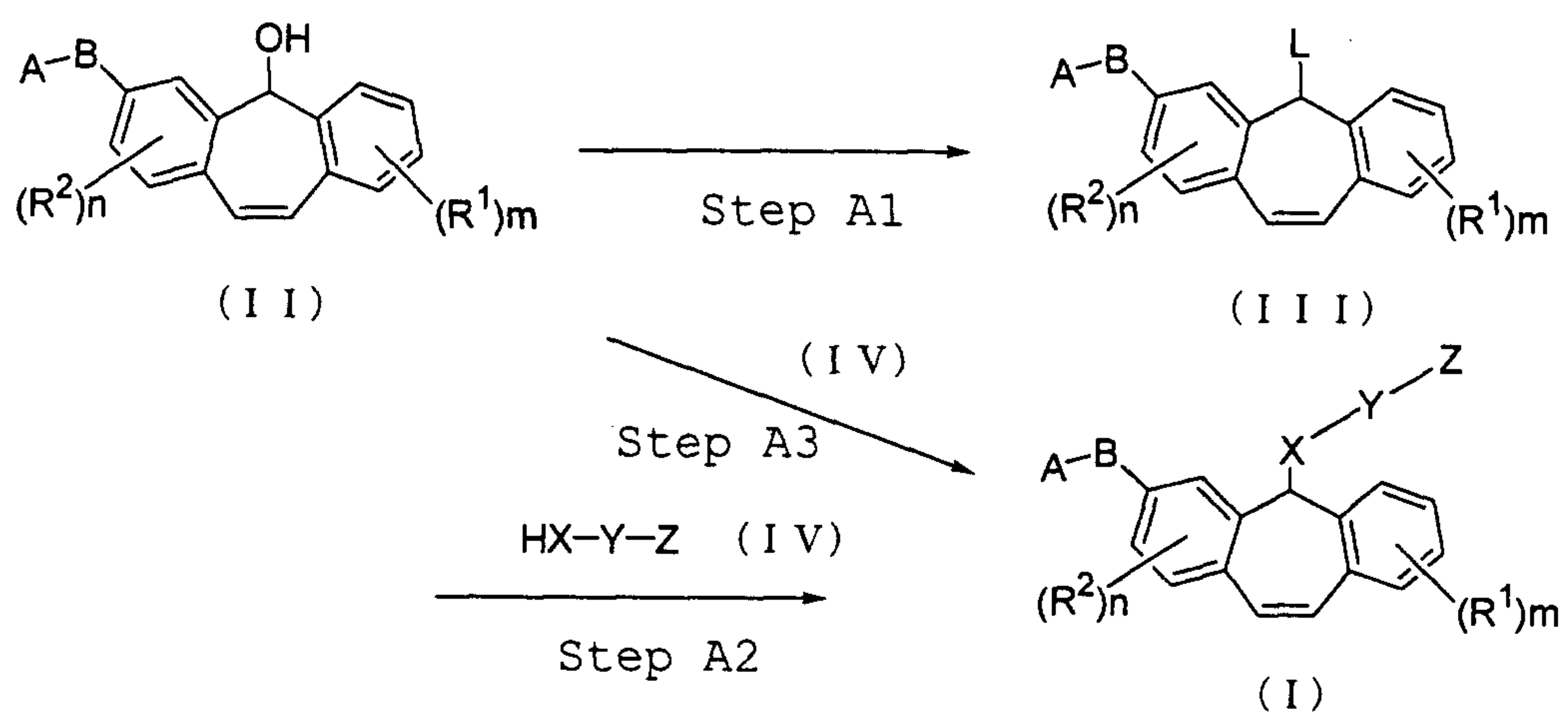
30 Compound No.1921; 3-{{[3-[(E)-2-(5,6,7,8-tetrahydroquinolin-2-yl)ethenyl]-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-yl]thio} propionic acid.

The compound represented by the formula (I) of the present invention can be produced by, for example,

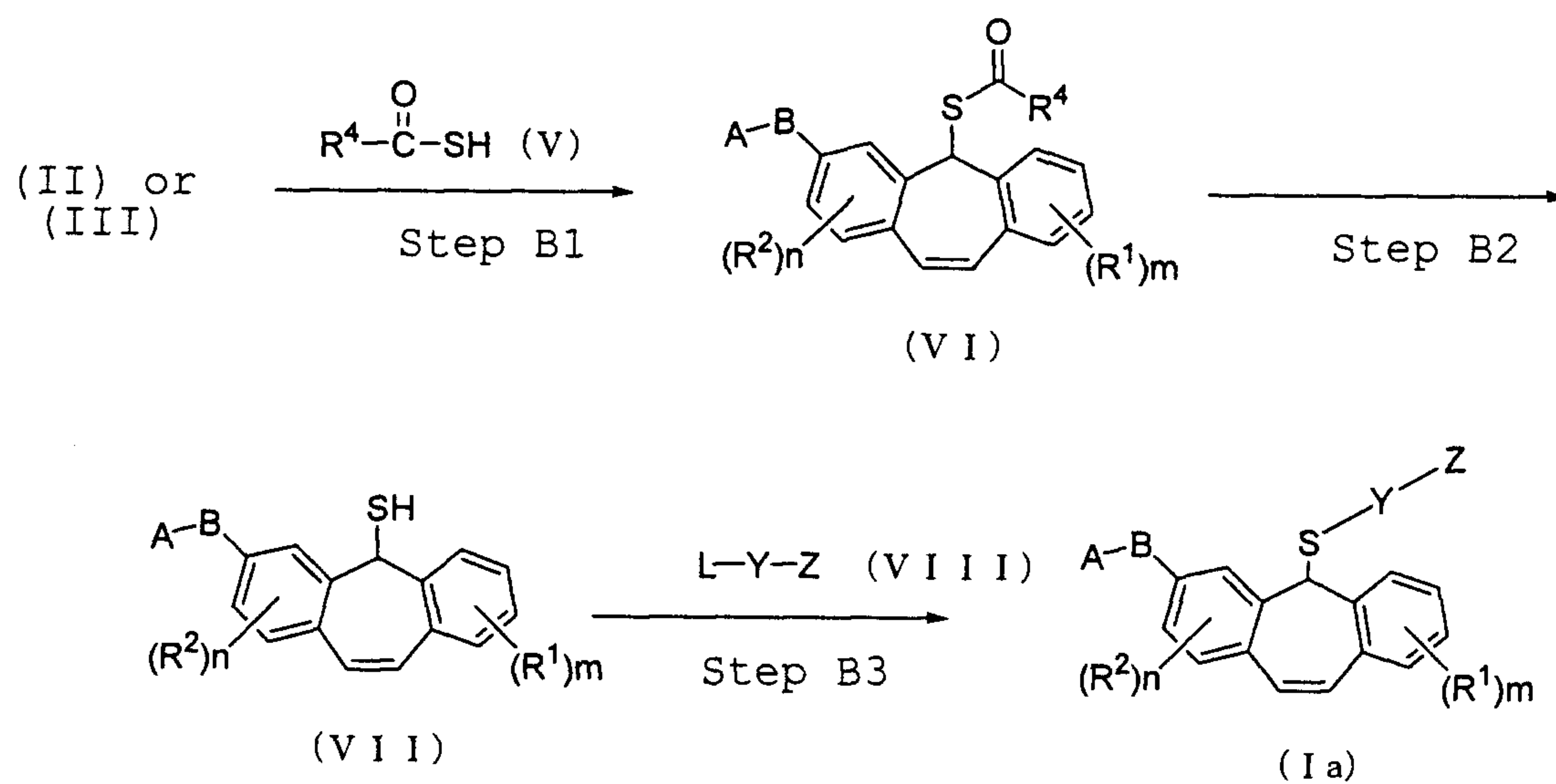
35 Preparation process A, B, C, D, E or L shown below.

100

Preparation process A

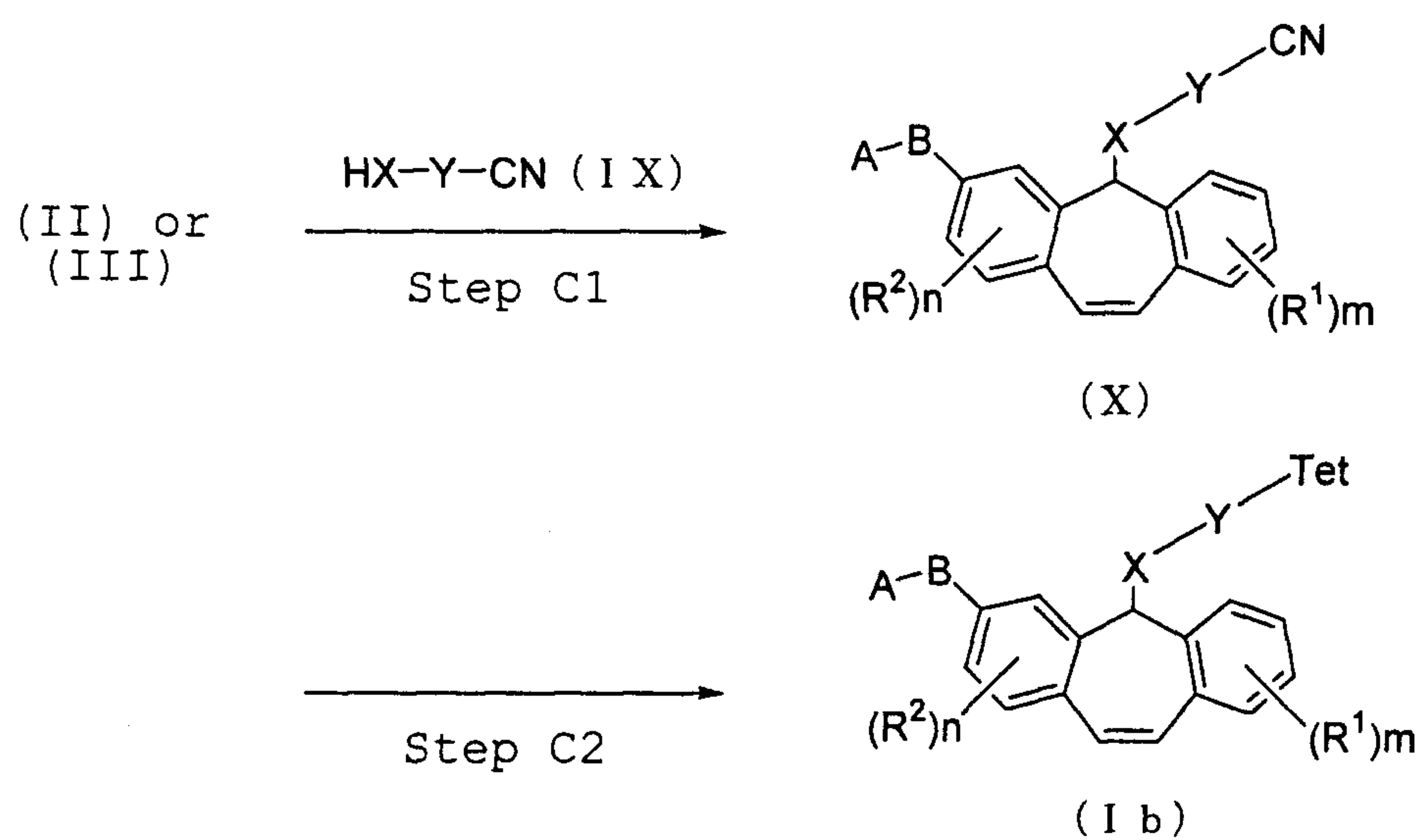


5 Preparation process B

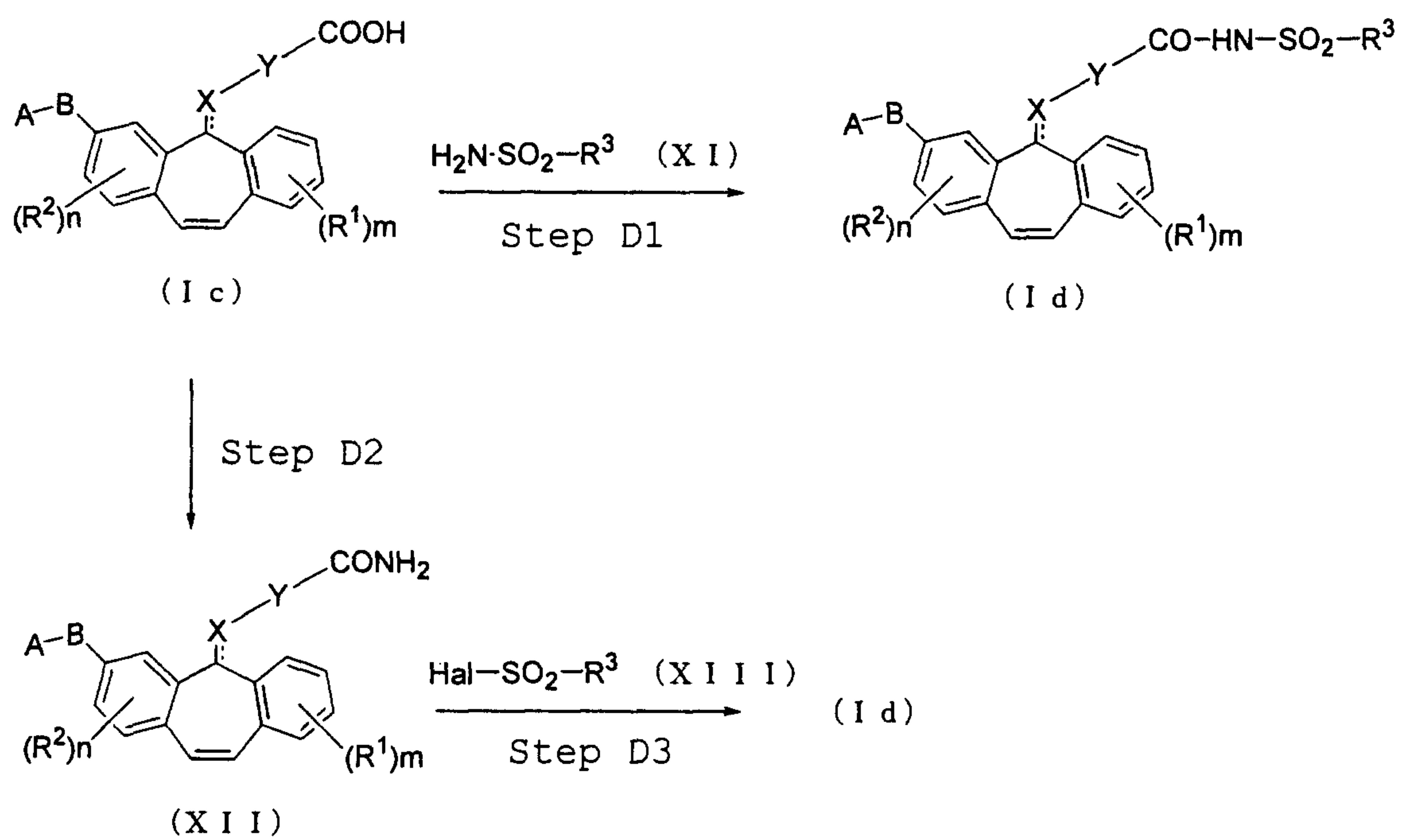


Preparation process C

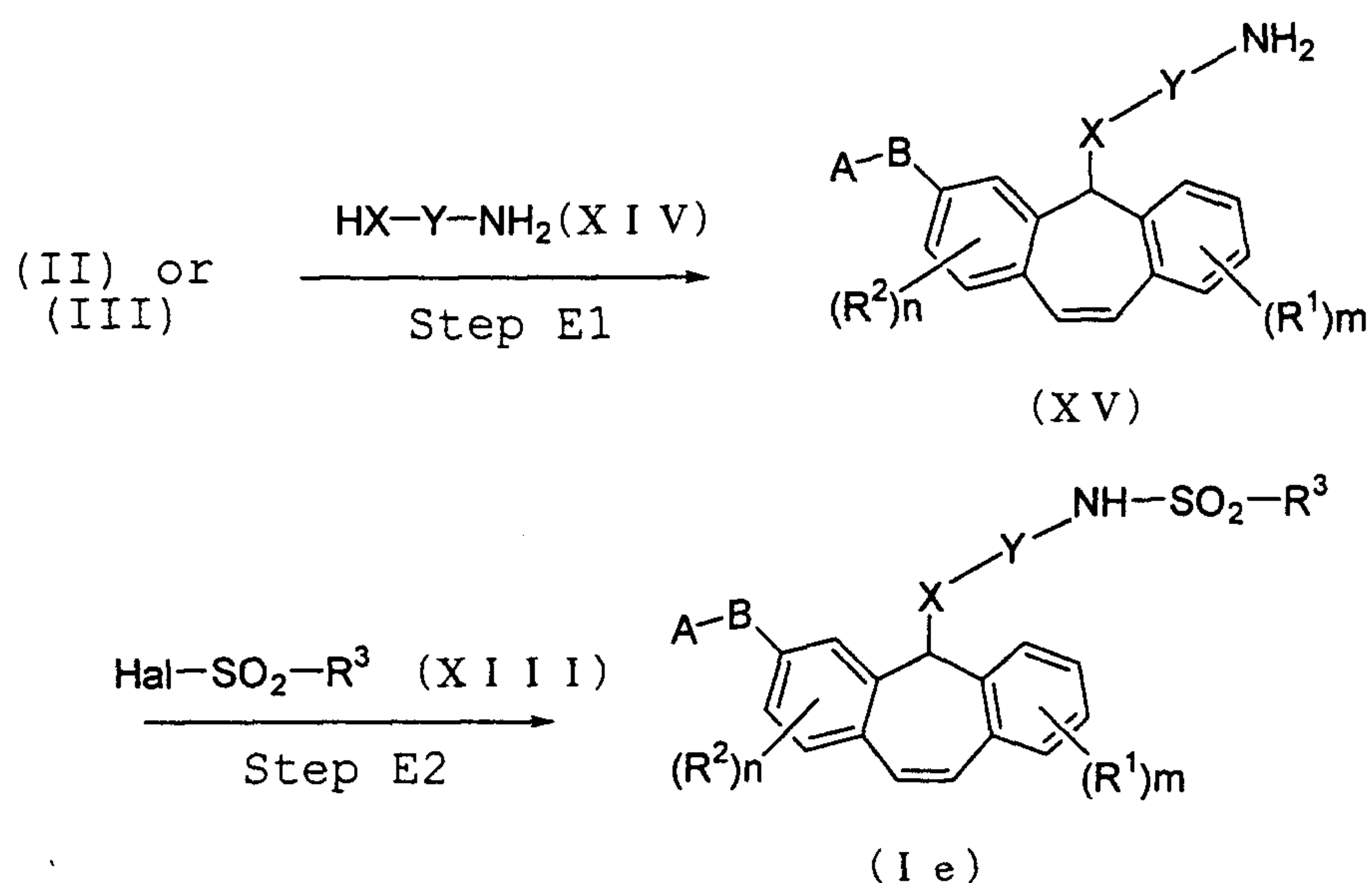
10



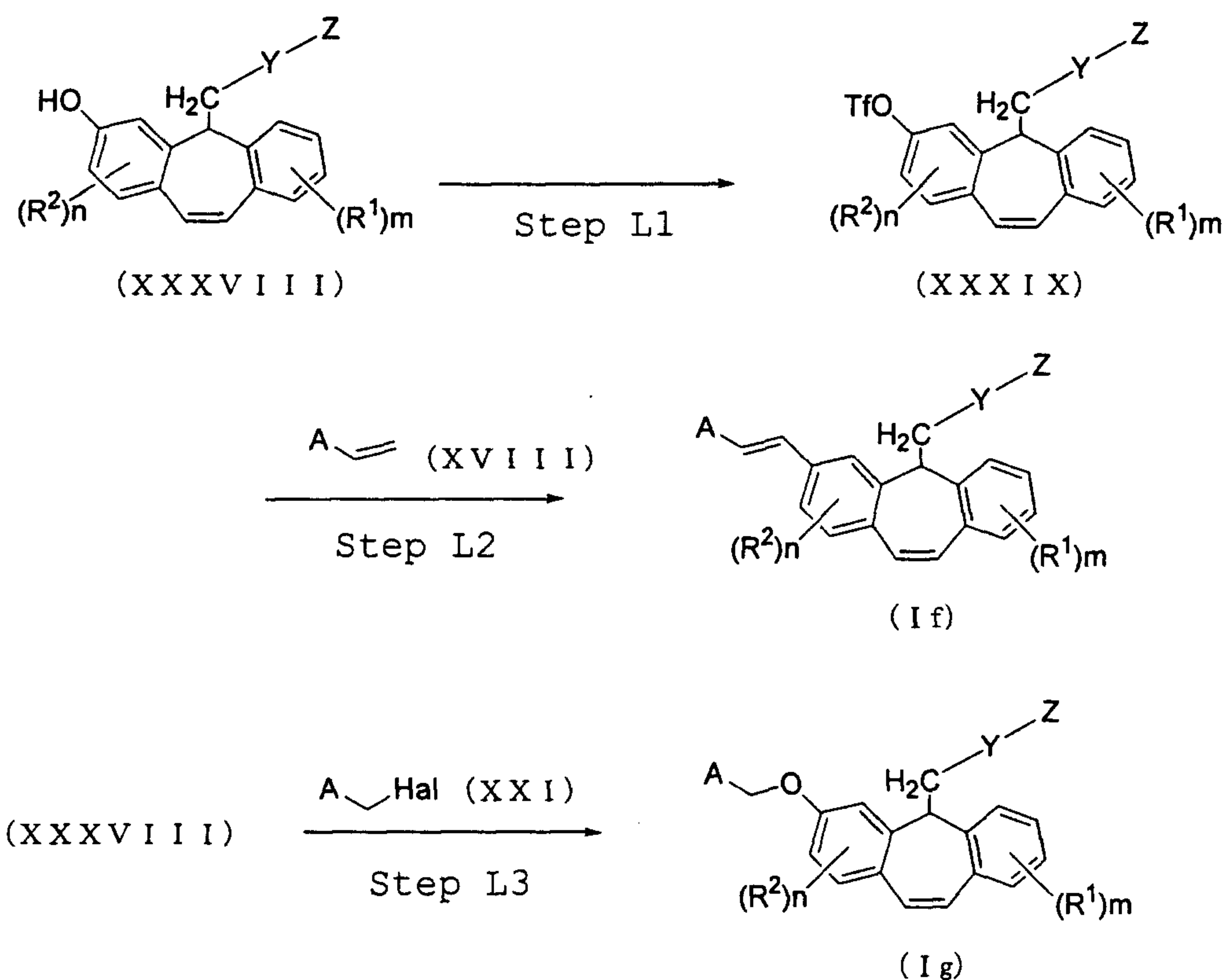
Preparation process D



Preparation process E



Preparation process L



5

In the above reaction formulae, R¹, R², R³, A, B, X, Y, Z, m and n have the same meanings as defined above, L represents a halogen atom, a C₁-C₄ alkylsulfonyloxy group, a fluoro C₁-C₄ alkylsulfonyloxy group or a phenylsulfonyloxy group which may be substituted (said substituent is a

10

C₁-C₄ alkyl group or a halogen atom), R⁴ represents a C₁-C₄ alkyl group or a phenyl group which may be substituted (said substituent is a C₁-C₄ alkyl group or a halogen atom), Tet represents a 1H-tetrazol-5-yl group, and Hal represents a halogen atom.

Preparation process A is a preparation process of Compound (I).

Step A1 of Preparation process A is a step of synthesizing Compound (III) by subjecting Compound (II) to halogenation or sulfonylation.

Halogenation of Compound (II) is carried out by reacting Compound (II) and a halogenating agent in a solvent or in the absence of a solvent (preferably in a solvent).

The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, halogenated hydrocarbons such as methylene chloride, chloroform, carbon tetrachloride, 1,2-dichloroethane, etc.; aromatic hydrocarbons such as benzene, toluene, etc.; or aliphatic hydrocarbons such as heptane, hexane, cyclohexane, etc., preferably halogenated hydrocarbons.

As the halogenating agent, there may be mentioned, for example, thionyl chloride, thionyl bromide, phosphorus oxychloride, phosphorus oxybromide, phosphorus pentachloride, etc., preferably thionyl chloride or phosphorus oxychloride. An amount of the halogenating agent to be used is usually 1 to 10-fold mol amount, preferably 1 to 2-fold mol amount based on the amount of Compound (II).

The reaction is carried out in the temperature range usually at -20 to 100°C, preferably at -10 to 30°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 5 minutes to 10 hours, preferably 10 minutes to 5 hours.

Sulfonylation of Compound (II) is carried out by

reacting Compound (II) and a sulfonylating agent in the presence of a base in a solvent.

The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, the same solvent as used in the above-mentioned halogenation reaction (for example, halogenated hydrocarbons, aromatic hydrocarbons or aliphatic hydrocarbons), etc., preferably halogenated hydrocarbons.

As the sulfonylating agent, there may be mentioned, for example, methanesulfonyl chloride, trifluoromethanesulfonyl chloride, methanesulfonic anhydride, trifluoromethanesulfonic anhydride, benzenesulfonyl chloride, toluenesulfonyl chloride, benzenesulfonyl bromide, toluenesulfonyl bromide, etc., preferably methanesulfonyl chloride, benzenesulfonyl chloride or toluenesulfonyl chloride. An amount of the sulfonylating agent to be used is usually 1 to 10-fold mol amount, preferably 1 to 3-fold mol amount based on the amount of Compound (II).

As the base, there may be mentioned, for example, amines such as triethylamine, tributylamine, diisopropylethylamine, pyridine, picoline, lutidine, 4-dimethylaminopyridine, etc., preferably triethylamine, diisopropylethylamine or pyridine. An amount of the base to be used is usually 1 to 10-fold mol amount, preferably 1 to 2-fold mol amount based on the amount of the sulfonylating agent.

The reaction is carried out in the temperature range usually at -10 to 100°C, preferably at 0 to 30°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 5 minutes to 10 hours, preferably 30 minutes to 5 hours.

Incidentally, Compound (III) can be separated and purified from the reaction mixture according to the usual method, but a crude product obtained by concentrating the reaction mixture can be used as such in the next step.

Step A2 is carried out by reacting Compound (III) and Compound (IV) in the presence or absence of a base in a solvent.

5 An amount of Compound (IV) to be used is usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount based on the amount of Compound (III).

10 The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, aprotic polar solvents such as N,N-dimethylformamide, N-methylpyrrolidone, dimethylsulfoxide, N,N-dimethylacetamide, hexamethylphosphoric acid triamide, etc.; halogenated hydrocarbons such as methylene chloride, chloroform, 1,2-dichloroethane, etc.; ketones such as acetone, methyl ethyl ketone, methyl isobutyl ketone, etc.; nitriles such as acetonitrile, etc.; esters such as ethyl acetate, etc.; or ethers such as diethyl ether, diisopropyl ether, tetrahydrofuran, dioxane, etc. or a mixed solvent of the above-mentioned solvents, 15 preferably aprotic polar solvents, ethers or a mixed solvent of the above-mentioned solvents.

20 As the base to be used, there may be mentioned, for example, alkali metal hydrides such as sodium hydride, lithium hydride, etc.; alkali metal amides such as sodium amide, etc.; amines such as triethylamine, tributylamine, diisopropylethylamine, pyridine, picoline, lutidine, 4-dimethylaminopyridine, etc.; or alkali metal carbonates such as sodium carbonate, potassium carbonate, sodium hydrogen carbonate, etc., preferably amines or alkali metal hydrides. An amount of the base to be used may vary 25 depending on the kinds of the starting compound(s), and it is usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount based on the amount of Compound (IV). Incidentally, in the present reaction, the base may not 30 necessarily be used.

The reaction is carried out in the temperature range

usually at -50 to 150°C, preferably at -10 to 100°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 5 minutes to 10 hours, preferably 30 minutes to 5 hours.

5 Step A3 is a step of separately obtaining Compound (I), and it is suitably employed particularly when X is a sulfur atom. The present step is carried out by reacting Compound (II) and Compound (IV) in the presence of an acid catalyst in a solvent.

10 An amount of Compound (IV) to be used is usually 1 to 5-fold mol amount, preferably 1 to 2-fold mol amount based on the amount of Compound (II).

 The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, halogenated hydrocarbons such as methylene chloride, chloroform, 1,2-dichloroethane, etc.; alcohols such as methanol, ethanol, propanol, isopropanol, butanol, etc.; aprotic polar
15 solvents such as N,N-dimethylformamide, N-methylpyrrolidone, dimethylsulfoxide, N,N-dimethylacetamide, hexamethylphosphoric acid triamide, etc.; or ethers such as diethyl ether, diisopropyl ether, tetrahydrofuran, dioxane, etc., preferably halogenated hydrocarbons.

25 As the acid catalyst to be used, there may be mentioned, for example, mineral acids such as hydrochloric acid, sulfuric acid, phosphoric acid, etc.; organic acids such as methanesulfonic acid, trifluoroacetic acid, etc.; Lewis acids such as boron trifluoride-diethyl ether complex, zinc chloride, tin tetrachloride, aluminum chloride,
30 etc., preferably organic acids or boron trifluoride-diethyl ether complex. An amount of the catalyst to be used is usually 0.1 to 50-fold mol amount, preferably 1 to 10-fold mol amount based on the amount of Compound (II), and when
35 the organic acids are used, it can be used with a markedly excessive amount also as a solvent.

The reaction is carried out in the temperature range usually at -10 to 100°C, preferably at 0 to 30°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 5 minutes to 10 hours, preferably 10 minutes to 5 hours.

Incidentally, in Compound (I), a compound wherein Z is a carboxyl group (Compound Ic mentioned below) can be directly produced by using Compound (IV) wherein Z is a carboxyl group, or can be synthesized, by using Compound (IV) wherein Z is a protected carboxyl group (the protective group is preferably a C₁-C₄ alkyl group), by once leading to Compound (I) wherein Z is a protected carboxyl group, and then, hydrolyzing the protective group under acidic or alkaline conditions by the conventional manner.

Also, in Compound (I), a desired protective group can be introduced into Compound (Ic) wherein Z is a carboxyl group according to the conventionally known method (for example, see written by W.Greene and P.G.H.Wult, "Protective Group in Organic Synthesis" 2nd Ed., John Wiley & Sons, p. 224).

Preparation process B is a preparation process of Compound (Ia) wherein X is a sulfur atom in Compound (I).

A reaction of obtaining Compound (VI) from Compound (II) or Compound (III) and a thiocarboxylic acid (V) in Step B1 can be carried out in the same manner as in the method described in Step A2 or Step A3 in the above-mentioned Preparation process A except for using a thiocarboxylic acid (V) in place of Compound (IV).

In Step B2, Compound (VII) can be synthesized by hydrolyzing Compound (VI) under alkaline conditions according to the conventional manner.

Step B3 is carried out by reacting Compound (VII) and Compound (VIII) in the presence of a base in a solvent.

The present reaction can be carried out in the same manner as in the method described in the above-mentioned Step A2 except for using Compound (VII) in place of Compound (II),

and using Compound (VIII) in place of Compound (IV).

An amount of Compound (VIII) to be used is usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount based on the amount of Compound (VII).

5 Preparation process C is a preparation process of Compound (Ib) wherein Z is a 1H-tetrazole group in Compound (I), and in Step C1, a reaction of obtaining Compound (X) from Compound (II) or Compound (III) and Compound (IX) can be carried out in the same manner as in the method
10 described in the above-mentioned Process A except for using Compound (IX) in place of Compound (IV).

In Step C2, Compound (Ib) can be synthesized by reacting Compound (X) and an azide compound in a solvent.

As the azide compound to be used, there may be
15 mentioned, for example, alkali metal azides such as sodium azide, potassium azide, lithium azide, etc.; alkaline earth metal azides such as calcium azide, magnesium azide, etc.; or organic tin azides such as trimethyl tin azide, tributyl tin azide, triphenyl tin azide, etc. An amount of the
20 azide compound to be used, it is usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount based on the amount of Compound (X). In the reaction, the azide compound is used singly, or may be used in combination with, for example, Lewis acids such as aluminum chloride, stannic
25 chloride, zinc chloride, titanium chloride, boron trifluoride-diethyl ether complex, etc.; ammonium salts such as ammonium chloride, tetramethylammonium chloride, etc.; sulfonic acids such as methanesulfonic acid, ethanesulfonic acid, etc.; alkali metal chlorides such as lithium chloride, etc.; or amine salts such as triethylamine hydro-
30 chloride, etc.

The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and
35 there may be mentioned, for example, aprotic polar solvents such as N,N-dimethylformamide, dimethylsulfoxide, N-methyl-

pyrrolidone, N,N-dimethylacetamide, etc.; ethers such as tetrahydrofuran, dimethoxyethane, diethoxyethane, dioxane, etc.; aromatic hydrocarbons such as benzene, toluene, xylene, etc.; or aliphatic hydrocarbons such as hexane, petroleum ether, etc.

The reaction is carried out in the temperature range usually at 0 to 200°C, preferably 50 to 150°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 1 hour to 72 hours, preferably 3 hours to 48 hours.

Preparation process D is a preparation process of Compound (Id) wherein Z is a $-\text{CO}-\text{NH}-\text{SO}_2-\text{R}^3$ group in Compound (I). Step D1 comprises a method of reacting Compound (Ic) and Compound (XI) in the presence of a condensing agent (Step D1a), or once leading Compound (Ic) to its reactive derivative (Step D1b), and then, reacting the reactive derivative and Compound (XI) in the presence of a base (Step D1c).

As the condensing agent to be used in Step D1a, there may be mentioned, for example, N,N'-dicyclohexylcarbodiimide (DCC), 1-ethyl-3-(3-dimethylaminopropyl)carbodiimide hydrochloride (EDC), N,N'-carbonyldiimidazole (CDI), diphenylphosphoric acid azide, hexafluorophosphoric acid benzotriazol-1-yloxy-tris-(dimethylamino)phosphonium (BOP), hexafluorophosphoric acid benzotriazol-1-yloxy-tris-pyrrolidinophosphonium (PyBOP), hexafluorophosphoric acid 2-(1H-benzotriazol-1-yl)-1,1,3,3-tetramethyluronium (HBTU), tetrafluoroboric acid 2-(1H-benzotriazol-1-yl)-1,1,3,3-tetramethyluronium (TBTU), etc., preferably DCC or EDC. An amount of the condensing agent to be used is usually 1 to 5-fold mol amount, preferably 1 to 3-fold mol amount based on the amount of Compound (Ic).

The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, ethers such as diethyl

ether, tetrahydrofuran, dimethoxyethane, diethoxyethane, dioxane, etc.; nitriles such as acetonitrile, etc.; aprotic polar solvents such as N,N-dimethylformamide, dimethylsulfoxide, N-methylpyrrolidone, N,N-dimethylacetamide, etc.; or
5 halogenated hydrocarbons such as methylene chloride, chloroform, 1,2-dichloroethane, etc., and these can be used singly or as a mixed solvent.

The reaction is carried out in the temperature range usually at -20 to 100°C, preferably at 0 to 50°C. A reaction
10 time may vary depending on the reaction temperature and the like, and it is usually 30 minutes to 24 hours, preferably 1 hour to 10 hours.

As the reactive derivative of Compound (Ic) in Step D1b, there may be mentioned, for example, acid halide
15 derivatives of Compound (Ic) such as acid bromide compound or acid chloride compound of Compound (Ic); or reactive amide derivatives such as Compound (Ic) and imidazol, 3,5-dimethylpyrazol or triazole and Compound (Ic), etc., preferably acid halide derivative.

20 The acid halide compound of Compound (Ic) can be produced according to the conventional manner, and for example, it can be synthesized by reacting Compound (Ic) with a halogenating agent (for example, thionyl chloride, thionyl bromide, phosphorus pentachloride, etc.) in a
25 solvent (for example, halogenated hydrocarbons such as methylene chloride, chloroform, 1,2-dichloroethane, etc.).

Also, reactive amide derivatives of Compound (Ic) can be produced according to the conventional manner, and for example, in the case of an imidazolidated compound of
30 Compound (Ic), it can be synthesized by reacting Compound (Ic) with 1,1'-carbonyldiimidazole in a solvent (for example, ethers such as diethyl ether, tetrahydrofuran, dimethoxyethane, diethoxyethane, dioxane, etc.; nitriles such as acetonitrile, etc.; aprotic polar solvents such as
35 N,N-dimethylformamide, dimethylsulfoxide, N-methylpyrrolidone, N,N-dimethylacetamide, etc.; or halogenated hydro-

carbons such as methylene chloride, chloroform, 1,2-dichloroethane, etc.).

The reactive derivatives of Compound (Ic) can be used as such in the next Step D1c without separation after
5 formation thereof.

Amounts of the reactive derivative of Compound (Ic) in Step D1c and Compound (XI) to be used in the reaction with Compound (XI) are each usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount based on the
10 amount of Compound (Ic).

As the base to be used, there may be mentioned, for example, amines such as triethylamine, tributylamine, diisopropylethylamine, pyridine, picoline, lutidine, 4-dimethylaminopyridine, 1,8-diazabicyclo[5.4.0]-7-undecene,
15 1,5-diazabicyclo[4.3.0]-5-nonene, etc., preferably triethylamine, tributylamine or diisopropylethylamine. An amount of the base to be used is usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount based on the amount of Compound (Ic).

20 The reaction is carried out in the temperature range usually at 0 to 150°C, preferably at 10 to 100°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 5 minutes to 48 hours, preferably 30 minutes to 24 hours.

25 Also, Compound (Id) can be produced by a method which is through Compound (XII).

Step D2 is a step of obtaining Compound (XII) by amidating the carboxyl group of Compound (Ic), and can be carried out by optionally employing a method among the
30 conventionally known methods. For example, Compound (XII) can be easily produced by reacting the above-mentioned reactive derivative of Compound (Ic) and ammonia.

Step D3 is carried out by reacting Compound (XII) and Compound (XIII) in the presence of a base in a solvent.

35 An amount of Compound (XIII) to be used is usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount

based on the amount of Compound (XII).

As the solvent and the base to be used, the same as those used in the above-mentioned Step D1 can be used, and the reaction can be carried out under the same conditions as those of Step D1.

Preparation process E is a preparation process of Compound (Ie) wherein Z is a $-\text{NH}-\text{SO}_2-\text{R}^3$ group in Compound (I).

In Step E1, the reaction of obtaining Compound (XV) from Compound (II) or Compound (III) and Compound (XIV) can be carried out by the same method as those mentioned in the above-mentioned method A except for using Compound (XIV) in place of Compound (IV).

Step E2 is carried out by reacting Compound (XV) and Compound (XIII) in the presence of a base in a solvent.

An amount of Compound (XIII) to be used is usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount based on the amount of Compound (XV).

As the solvent to be used, the same solvent as those mentioned in the above-mentioned Step D1 may be mentioned, and preferably halogenated hydrocarbons or aprotic polar solvents.

As the base to be used, the same base as those mentioned in the above-mentioned Step D1 may be mentioned, and an amount of the base to be used is usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount based on the amount of Compound (XIV). Also, in the present reaction, the base may be used in an extremely excessive amount also as a solvent.

The reaction is carried out in the temperature range usually at -20 to 100°C , preferably at 0 to 50°C . A reaction time may vary depending on the reaction temperature and the like, and it is usually 5 minutes to 10 hours, preferably 30 minutes to 5 hours.

Preparation process L is a preparation process of Compound (If) wherein X is a methylene group and B is

-CH=CH- in Compound (I), and Compound (Ig) wherein X is a methylene group and B is -CH₂O- in Compound (I).

Step L1 is carried out in the same manner as in Step F1 mentioned below except for using Compound (XXXVIII) in place of Compound (XVI), and Step L2 is carried out in the same manner as in Step F2 mentioned below except for using Compound (XXXIX) in place of Compound (XVII).

Step L3 is carried out in the same manner as in Step G1 of preparation process G mentioned below except for using Compound (XXXVIII) in place of Compound (XVI).

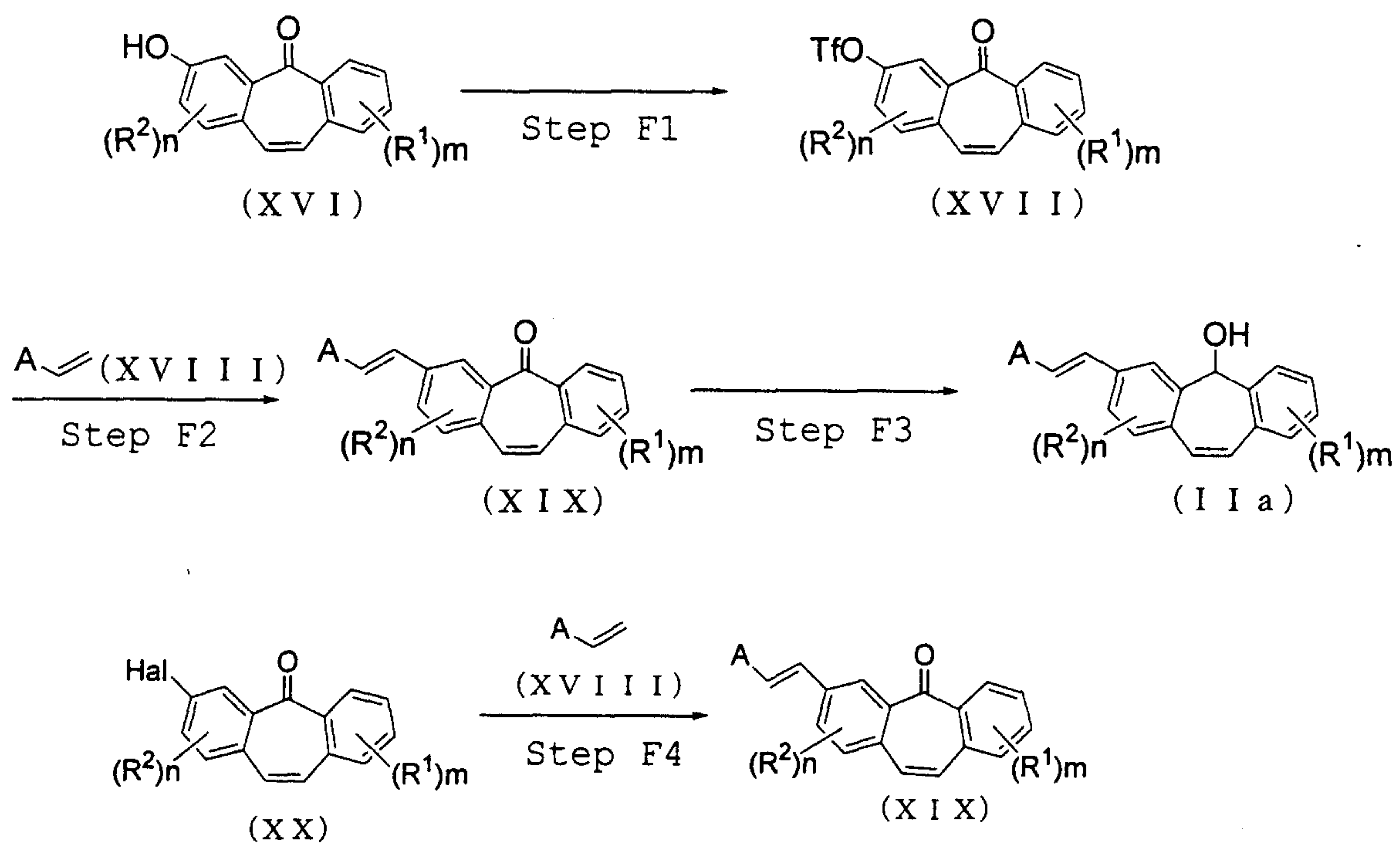
In the above-mentioned respective reactions, the formed objective Compound can be collected from the reaction mixture according to the conventional manner. For example, when insoluble materials are present, the objective compound can be obtained by, after removing the insoluble materials by filtration, removing the solvent, or by removing the solvent under reduced pressure, adding water to the residue, extracting the mixture with a water-immiscible organic solvent such as ethyl acetate, etc., if necessary, after drying over anhydrous sodium sulfate, etc., removing the solvent, and further, if necessary, by purifying with the conventional manner, for example, recrystallization, column chromatography, etc.

Also, the compound of the formula (I) of the present invention can be converted into a pharmaceutically acceptable salt by treating with an acid or a base according to the conventional manner. For example, in an inert solvent (preferably ethers such as diethyl ether, tetrahydrofuran, dimethoxyethane, diethoxyethane, dioxane, etc.; alcohols such as methanol, ethanol, propanol, isopropanol, butanol, etc.; halogenated hydrocarbons such as methylene chloride, chloroform, etc.; or water), it is reacted with a desired acid or a base, and the solvent is removed, or else, precipitated crystals are collected by filtration to obtain a desired salt. Also, it can be directly separated as a salt from the reaction mixture in the final reaction step.

Moreover, in the compound of the formula (I), there exist optical isomer(s) (including diastereomer) due to an asymmetric carbon(s) and/or geometric (E,Z) isomers due to an unsaturated carbon. These respective isomers can be
5 separated by treating the corresponding racemic isomers or geometric isomer mixture by usual optical resolution methods (fractional recrystallization method, optical resolution column chromatography method or diastereomer method, etc.) or separation methods (recrystallization
10 method, column chromatography method, etc.). For example, when optical isomers are to be separated, Compound (I) which is racemic mixture is reacted with an optically active sulfonic acid compound ((S) or (R)-camphor-10-sulfonic acid, etc.), to obtain one of the diastereomer
15 salts, and if necessary, further subjecting to purification, the resulting diastereomer salt is resolved according to the conventional manner to obtain an optical isomer. Also, when the above reaction is carried out by using the starting compound which has been subjected to optical
20 resolution or separation, a desired optical isomer or geometric isomer can be obtained.

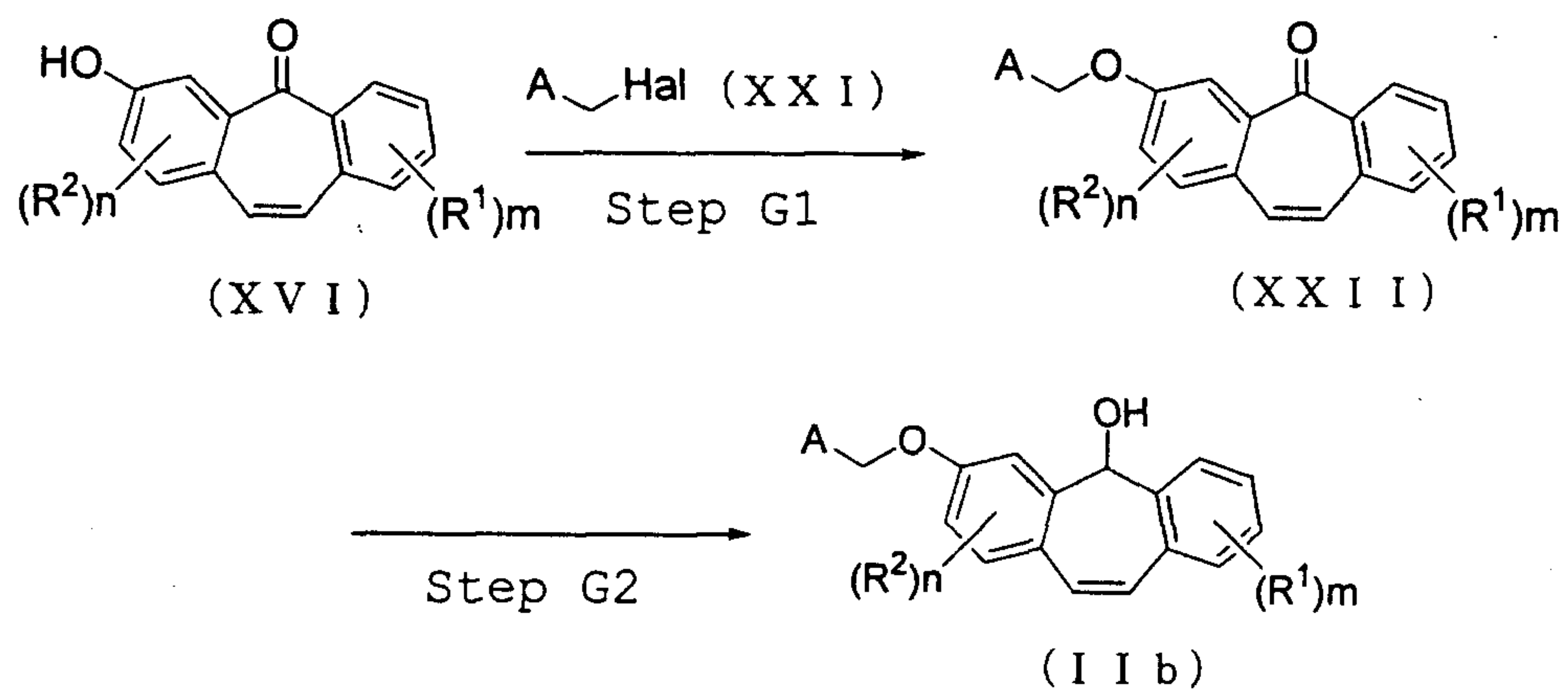
Compound (IV), (V), (VIII), (IX), (XI), (XIII), (XIV) and other sub-starting materials which are used as starting materials in the above-mentioned Preparation
25 process A, B, C, D, E or L are each well known compounds or can be easily produced according to the conventionally know method. Also, Compound (II) and (XXXVIII) can be produced by using either one of or in combination with two or more of Preparation processes F, G, H, I, J, K, N, O and P shown
30 below optionally.

Preparation process F



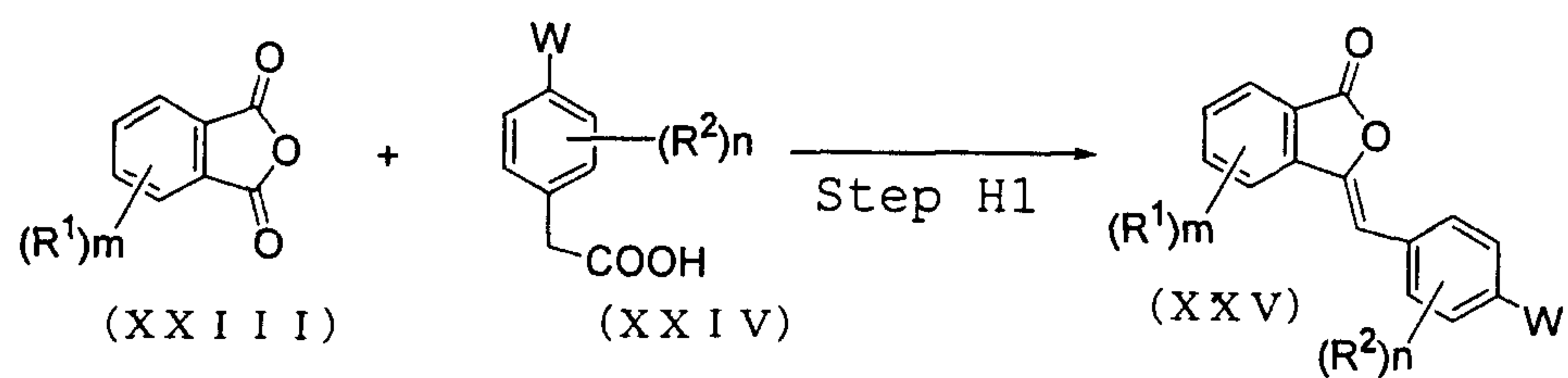
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Preparation process G

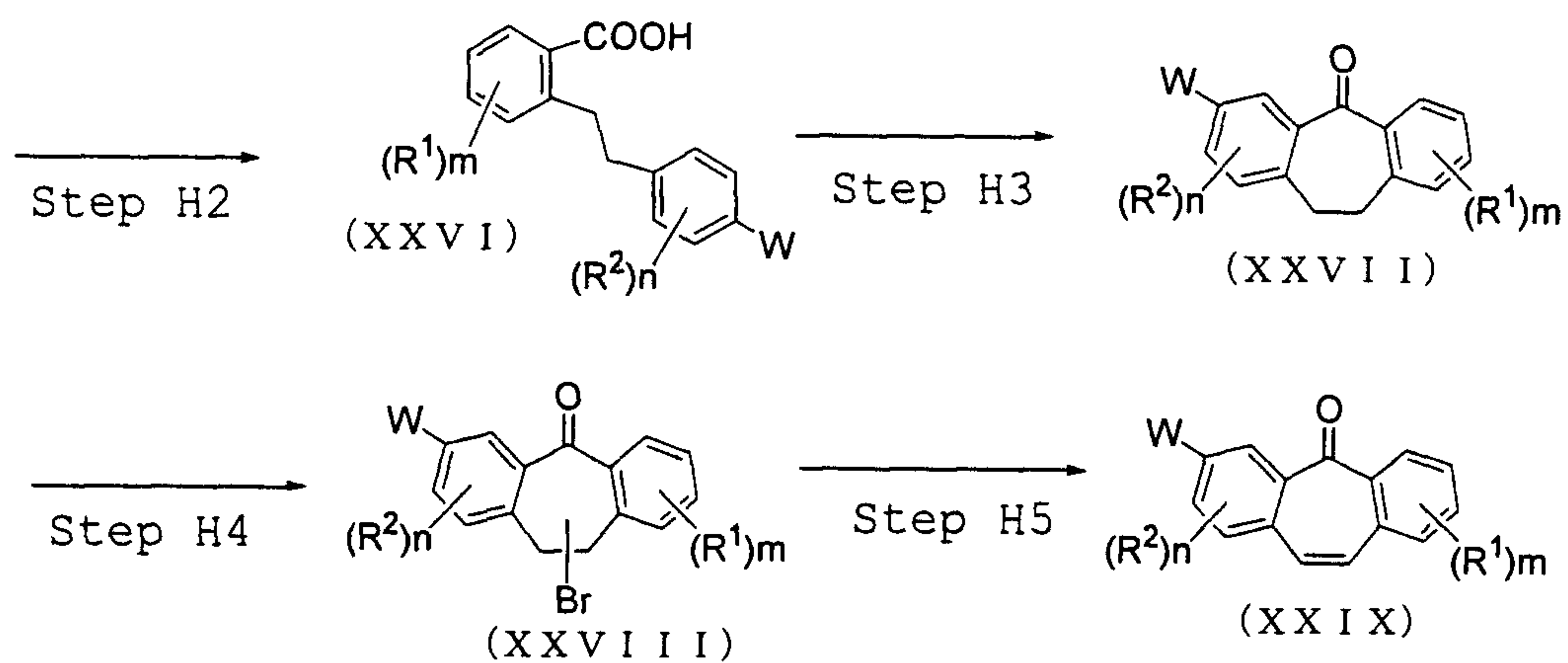


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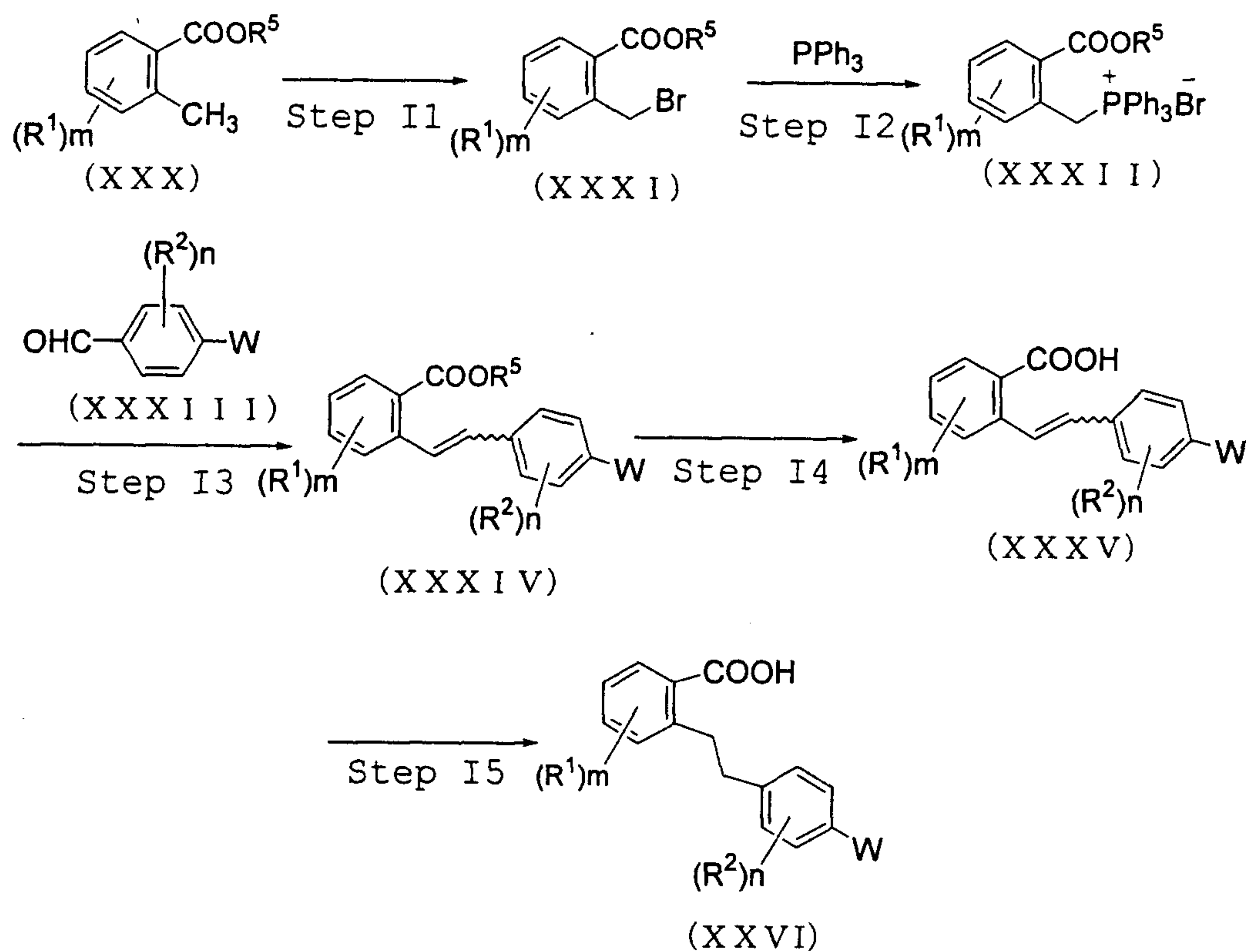
Preparation process H



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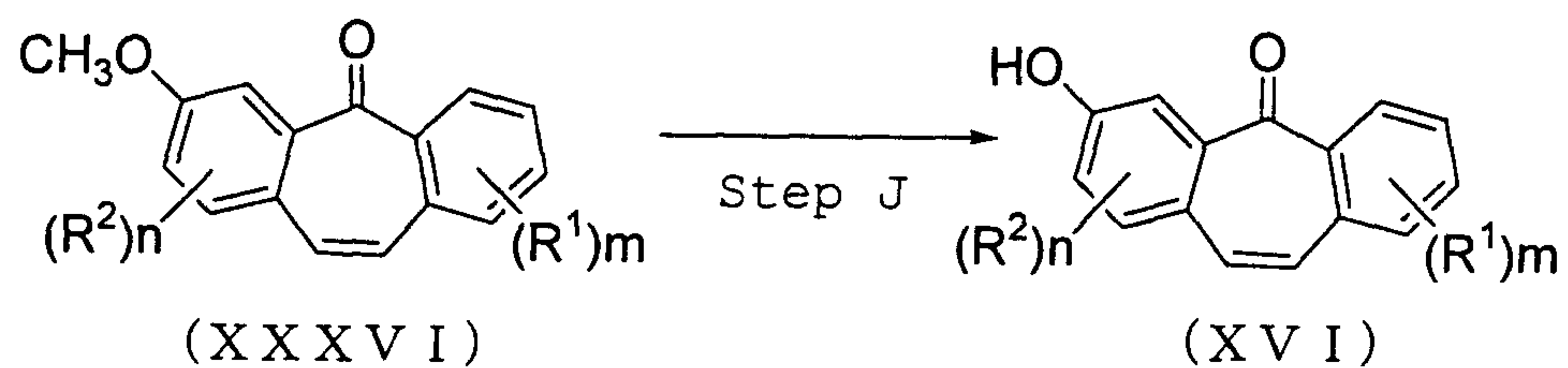


Preparation process I



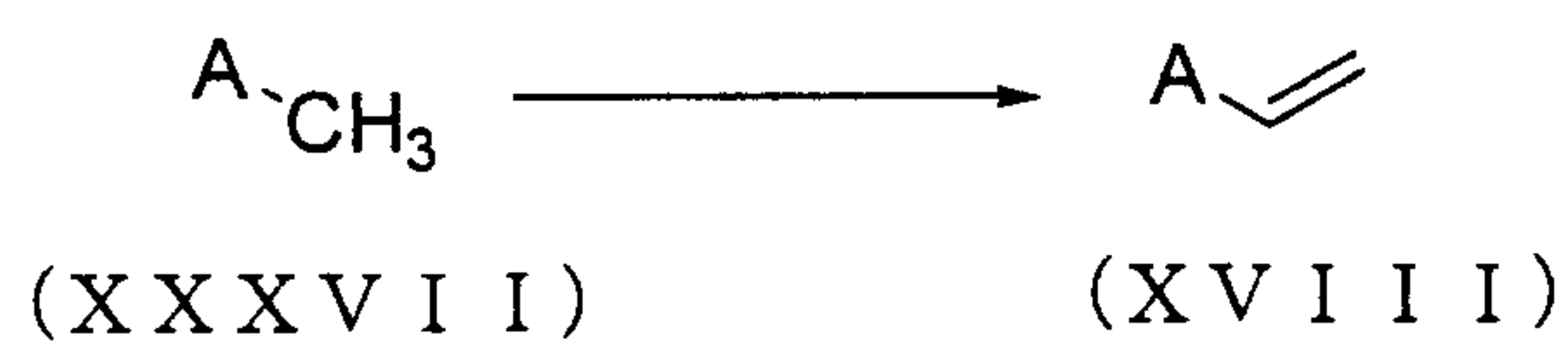
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Preparation process J



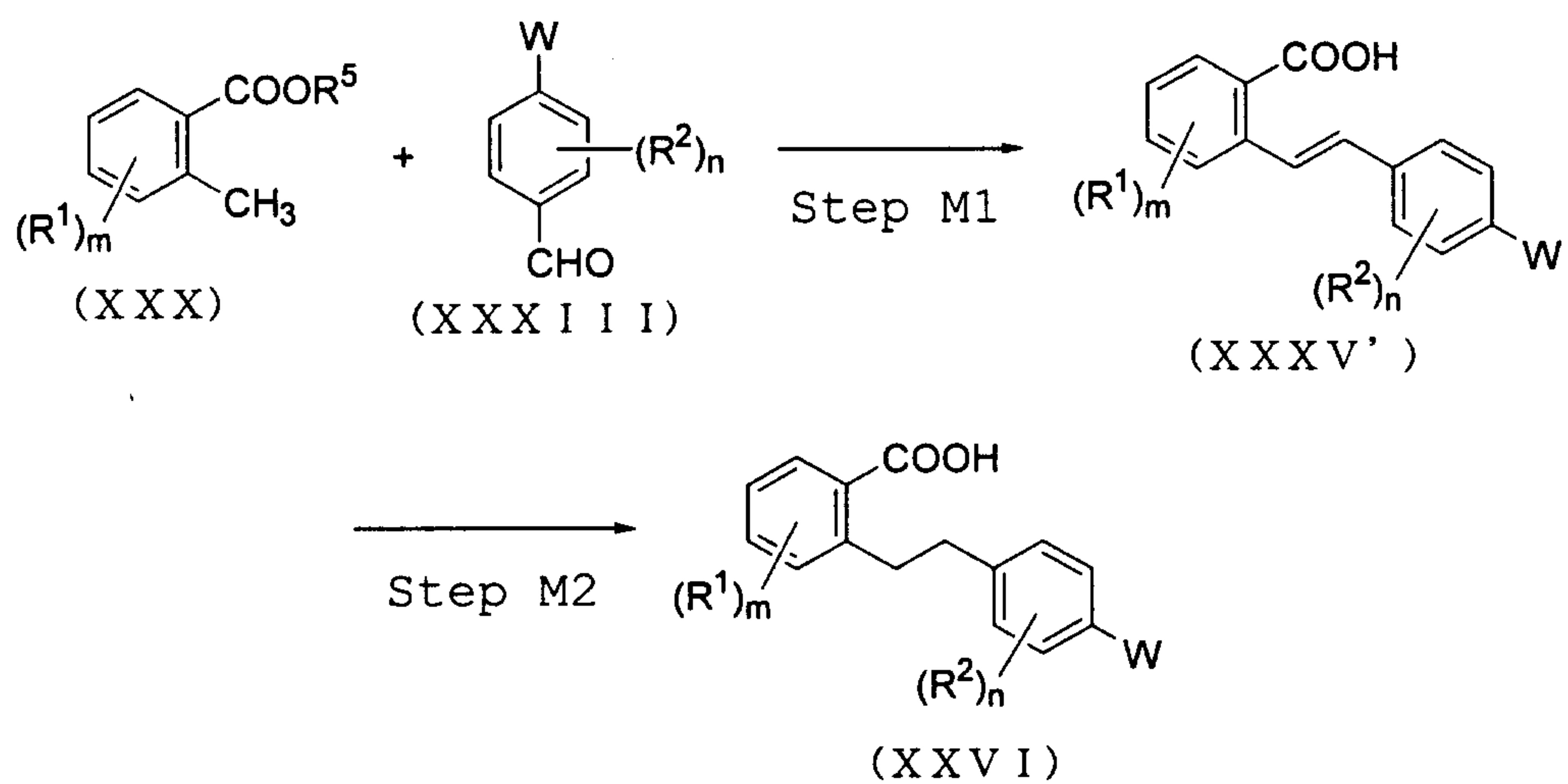
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Preparation process K



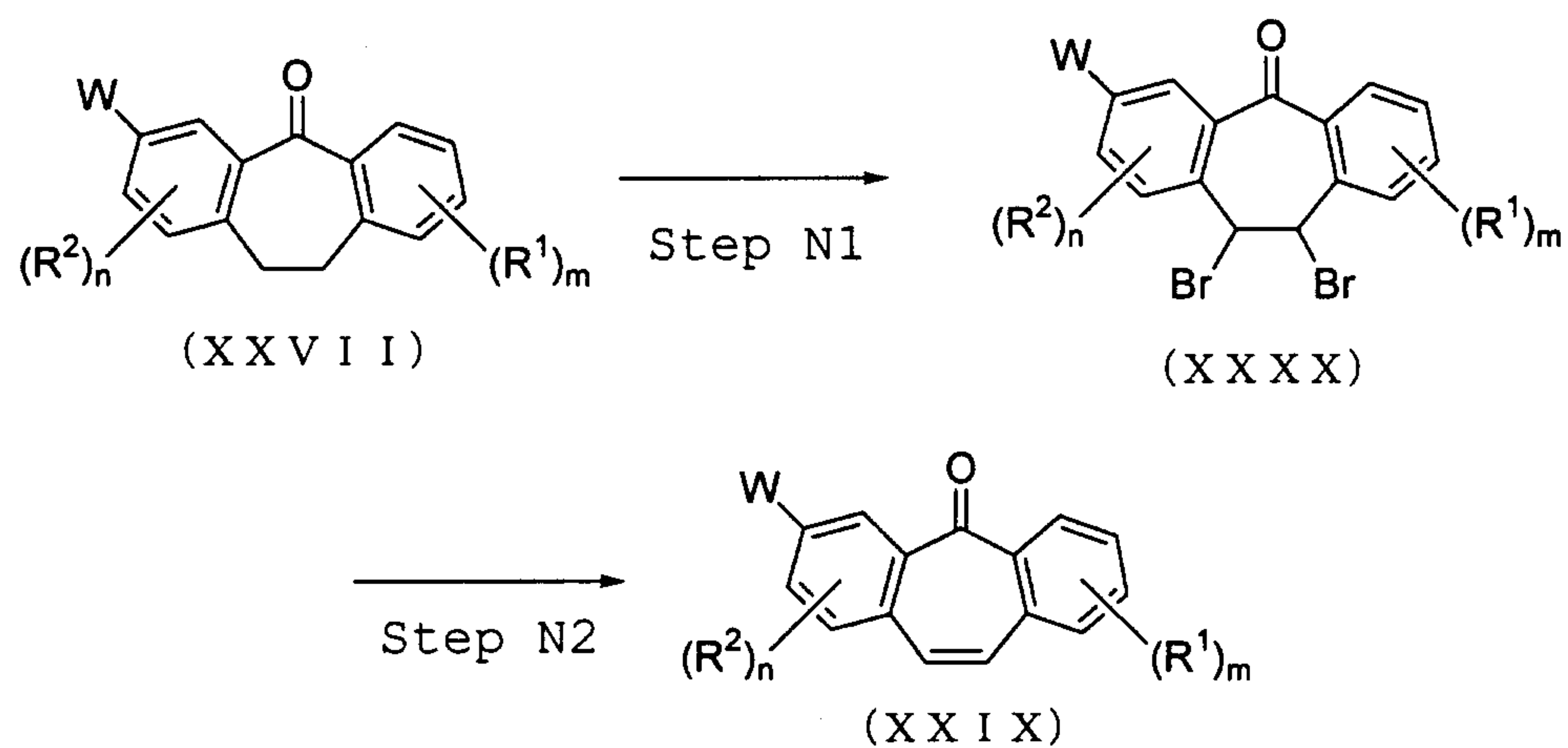
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Preparation process M

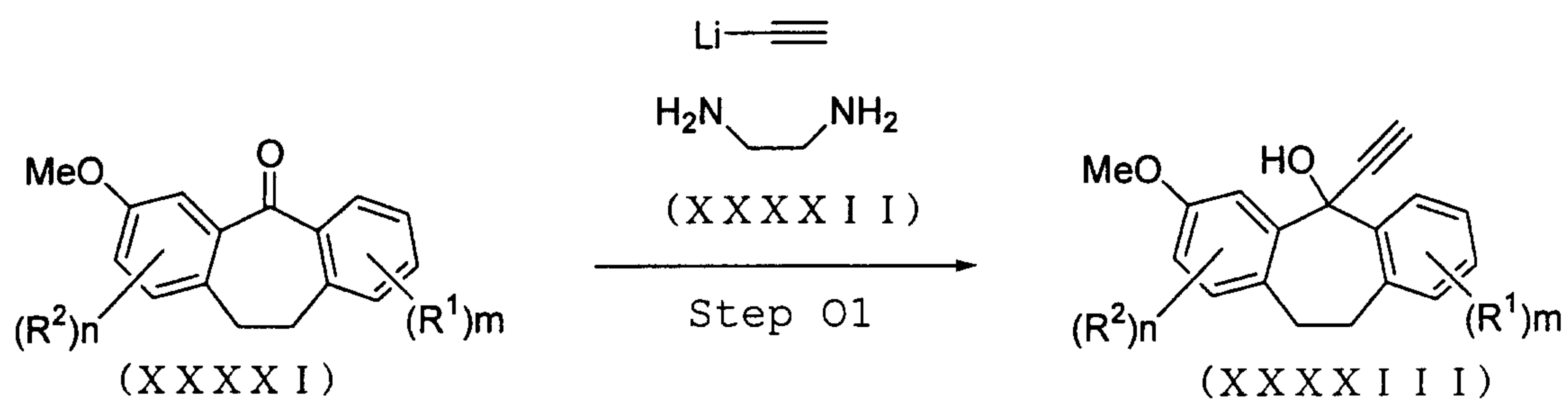


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Preparation process N

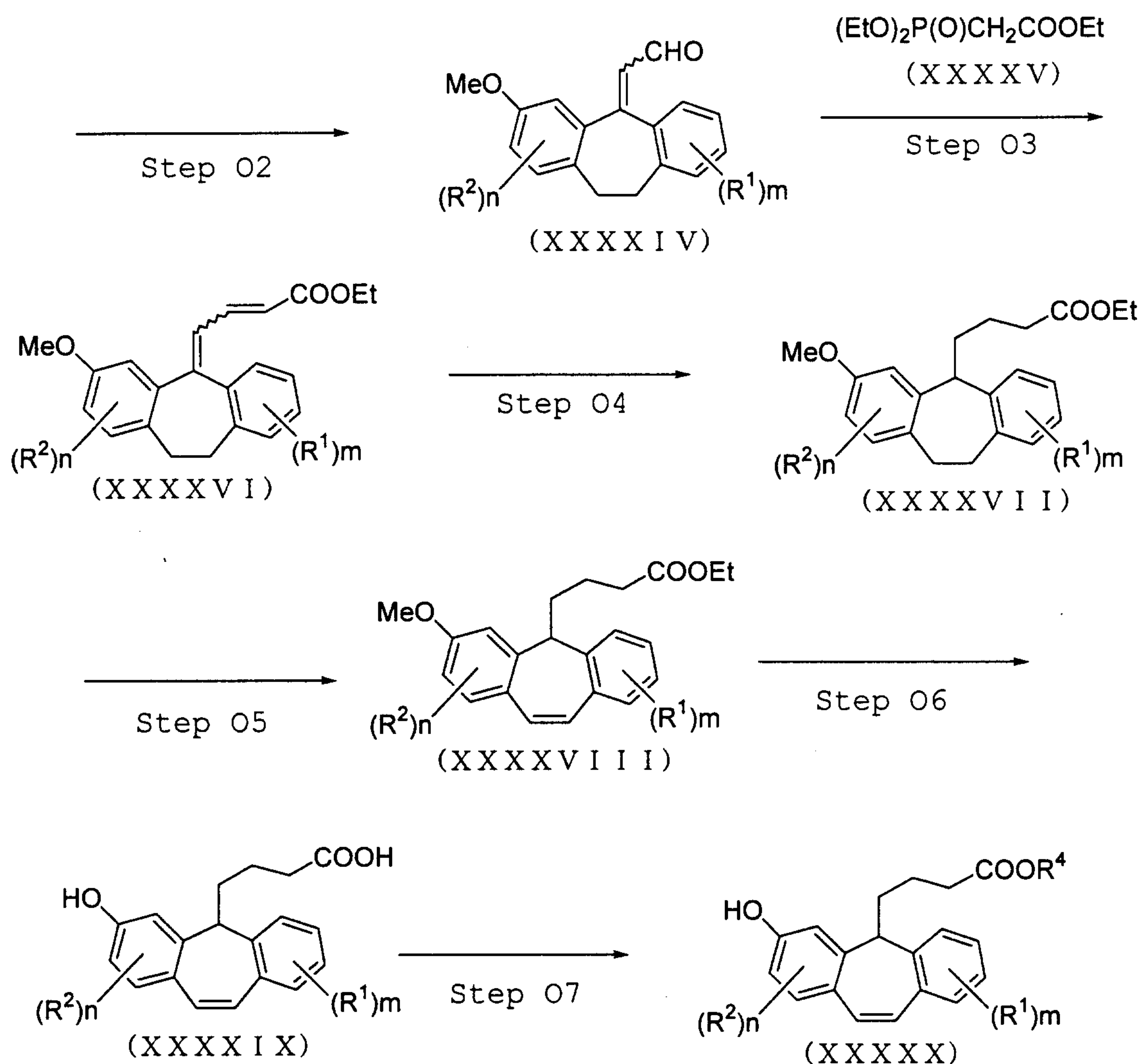


Preparation process O

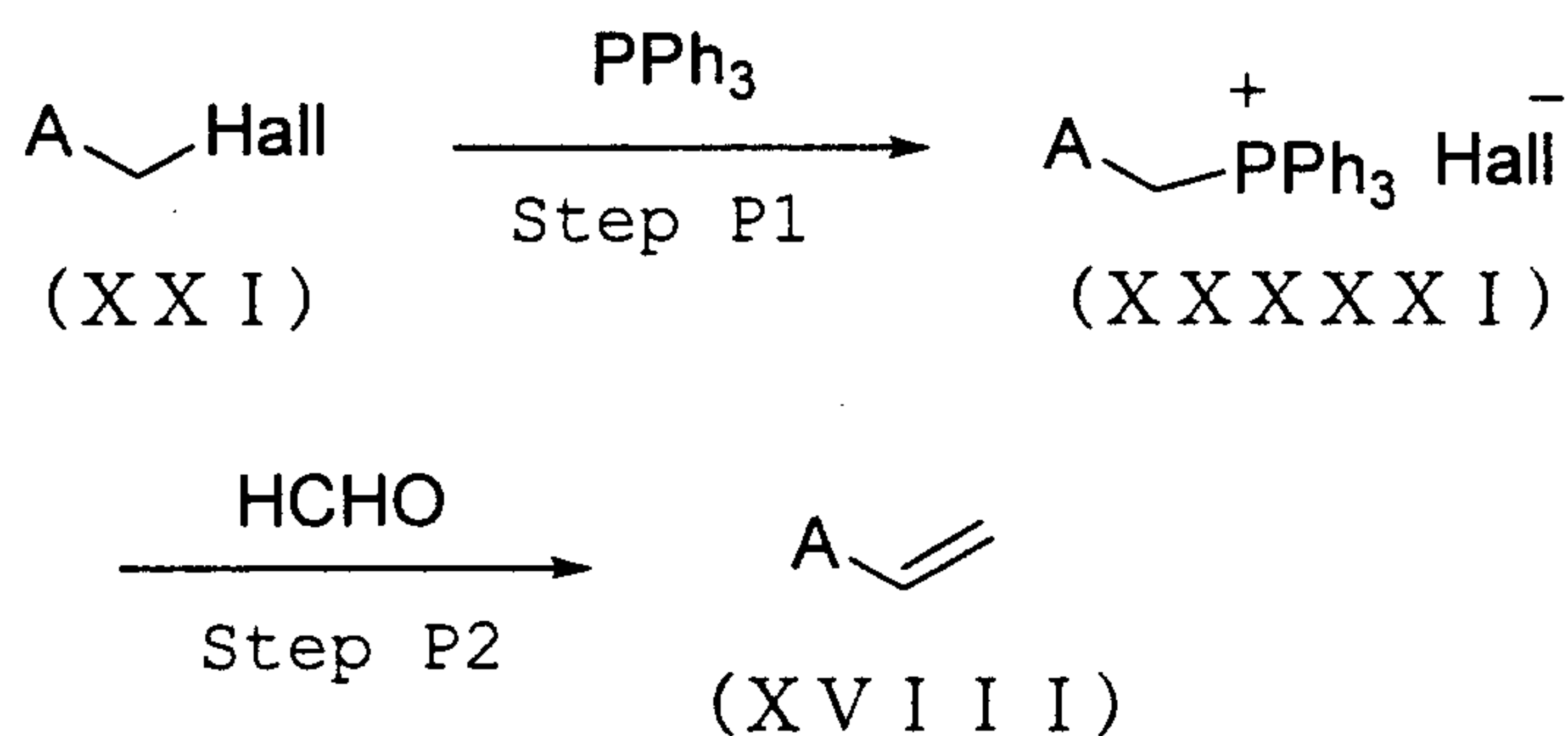


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Preparation process P



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wherein R^1 , R^2 , A, Hal, m and n have the same meanings as defined above, Tf represents a trifluoromethanesulfonyl group, W represents a methoxy group or a halogen atom (preferably a bromine atom or an iodine atom), and R^5 represents a C_1 - C_4 alkyl group.

10

Preparation process F is a preparation process of Compound (IIa).

Step F1 is a step of synthesizing Compound (XVII) by trifrating Compound (XVI).

5 Trifration of Compound (XVI) is carried out by reacting Compound (XVI) and a trifrating agent in the presence of a base in a solvent.

10 The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, halogenated hydrocarbons such as methylene chloride, chloroform, carbon tetrachloride, 1,2-dichloroethane, etc.; aromatic hydrocarbons such as benzene, toluene, etc.; or aliphatic
15 hydrocarbons such as heptane, hexane, cyclohexane, etc., preferably halogenated hydrocarbons.

20 As the trifrating agent, there may be mentioned, for example, trifluoromethanesulfonyl chloride or trifluoromethanesulfonic anhydride, preferably trifluoromethanesulfonic anhydride. An amount of the trifrating agent to be used is usually 1 to 10-fold mol amount, preferably 1 to 2-fold mol amount based on the amount of Compound (XVI).

25 As the base, there may be mentioned, for example, amines such as triethylamine, tributylamine, diisopropylethylamine, pyridine, picoline, lutidine, 4-dimethylaminopyridine, etc., preferably triethylamine, diisopropylethylamine or pyridine. An amount of the base to be used is usually 1 to 10-fold mol amount, preferably 1 to 2-fold mol amount based on the amount of the trifrating agent.

30 The reaction is carried out in the temperature range usually at -20 to 100°C, preferably at -10 to 30°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 5 minutes to 10 hours, preferably 30 minutes to 5 hours.

35 Step F2 is carried out by reacting Compound (XVII) and Compound (XVIII) in an inert gas atmosphere such as

nitrogen, helium or argon in the presence of a catalyst (palladium catalyst) and a base, in a solvent.

The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, polar solvents such as N,N-dimethylformamide, N,N-dimethylacetamide, N-methylpyrrolidone, dimethylsulfoxide, etc., or acetonitrile, etc., preferably N,N-dimethylformamide.

As the base, there may be mentioned, for example, amines such as triethylamine, tributylamine, diisopropylethylamine, pyridine, picoline, lutidine, 4-dimethylaminopyridine, etc., preferably triethylamine. An amount of the base to be used is usually 1 to 10-fold mol amount, preferably 1 to 2-fold mol amount based on the amount of Compound (XVII).

Also, in place of the amines, quaternary amines such as tetrabutyl ammonium chloride, tetrabutyl ammonium bromide, etc., and alkali metal carbonates such as potassium carbonate, sodium carbonate, sodium hydrogen carbonate, etc. may be used in combination.

As the palladium catalyst, there may be mentioned, for example, palladium acetate, palladium acetate-triphenylphosphine, dichlorobis(triphenylphosphine)palladium, tetrakis(triphenylphosphine)palladium, etc., preferably palladium acetate-triphenylphosphine or tetrakis(triphenylphosphine)palladium. An amount of the palladium catalyst to be used is usually 0.01 to 1-fold mol amount, preferably 0.01 to 0.3-fold mol amount based on the amount of Compound (XVII).

Also, lithium chloride or lithium bromide may be copresent in the reaction.

The reaction is carried out in the temperature range usually at 0°C to 200°C, preferably at 50°C to 150°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 30 minutes to 48

hours, preferably 1 hour to 24 hours.

Reduction from Compound (XIX) to Compound (IIa) in Step F3 can be carried out by using a reducing agent in a solvent.

5 As the reducing agent, there may be mentioned, for example, sodium borohydride, lithium borohydride, sodium cyanoborohydride, lithium aluminum hydride, etc., preferably lithium borohydride.

10 The solvent to be used may be mentioned, for example, alcohols such as methanol, ethanol, propanol, butanol, etc.; ethers such as tetrahydrofuran, dioxane, 1,2-dimethoxyethane, etc.; nitriles such as acetonitrile, etc.; amides such as N,N-dimethylformamide, N,N-dimethylacetamide, N-methylpyrrolidone, etc.; water or a mixed solvent
15 of the above-mentioned solvents, preferably tetrahydrofuran or a mixed solvent with tetrahydrofuran.

20 The reaction is carried out in the temperature range usually at -10 to 150°C, preferably at 0 to 100°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 10 minutes to 10 hours, preferably 30 minutes to 6 hours.

25 Step F4 is another process for producing Compound (XIX), and carried out in the similar conditions as those mentioned in the above-mentioned Step F2 by using Compound (XX) (preferably a bromine atom or an iodine atom as Hal) in place of Compound (XVII) as a starting compound.

Preparation process G is a preparation process of Compound (IIb).

30 Step G1 is carried out by reacting Compound (XVI) and Compound (XXI) in the presence of a base in a solvent, and carried out in the similar method as those mentioned in the above-mentioned Step A2 except for using Compound (XVI) in place of Compound (III), and using Compound (XXI) in place of Compound (IV).

35 Reduction from Compound (XXII) to Compound (IIb) in Step G2 is carried out in a solvent by using a reducing

agent. It is carried out in the similar method as those mentioned in the above-mentioned Step F3 except for using Compound (XXII) in place of Compound (XIX).

Preparation process H is a preparation process of
5 Compound (XXIX), and includes preparation processes of Compound (XX: W=a halogen atom) and a compound (XXXVI: W=methoxy group) mentioned below.

Preparation process H can be easily carried out by referring to the conventionally known method as disclosed
10 in, for example, *Helv. Chim. Acta*, 48, 1240 (1965) or USP 1,207,404 B.

Step H1 is carried out by, for example, reacting Compound (XXIII) and Compound (XXIV) in the presence of a base such as sodium acetate in the absence of a solvent or
15 using diphenyl ether as a solvent under heating (preferably at 100°C to 250°C) to carry out dehydration condensation.

A step of obtaining Compound (XXVI) by reducing Compound (XXV) in Step H2 is carried out by optionally employing a reducing method in which hydrogen is used in
20 the presence of a Raney-nickel catalyst (USP 1,207,404 B) or a reducing method in which hydroiodic acid/red phosphorus are used (*Helv. Chim. Acta*, 48, 1240 (1965)), and when W is a halogen atom, the latter reduction method is suitably used.

25 A step of obtaining Compound (XXVII) by subjecting Compound (XXVI) to dehydration and cyclization in Step H3 is carried out by reacting Compound (XXVI) in a solvent or in the absence of a solvent, and in the presence of a catalyst (a dehydrating agent).

30 The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, halogenated hydrocarbons such as methylene chloride, chloroform, carbon tetrachloride, 1,2-dichloroethane, etc.; nitrobenzene, and
35 carbon disulfide, preferably halogenated hydrocarbons.

As the catalyst to be used, there may be mentioned, for example, mineral acids such as sulfuric acid, phosphoric acid, polyphosphoric acid, etc.; acid anhydrides such as methanesulfonic anhydride, trifluoroacetic anhydride, etc.;
5 Lewis acids such as boron trifluoride-diethyl ether complex, aluminum chloride, zinc chloride, etc., preferably polyphosphoric acid, methanesulfonic anhydride, trifluoroacetic anhydride or boron trifluoride-diethyl ether complex. Also, a mixture of trifluoroacetic anhydride and
10 boron trifluoride-diethyl ether complex is suitably used. An amount of the catalyst to be used is usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount based on the amount of Compound (XXVI).

The reaction is carried out in the temperature range
15 usually at 0 to 150°C, preferably at 0 to 100°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 5 minutes to 10 hours, preferably 30 minutes to 5 hours.

For example, when polyphosphoric acid, etc. as the
20 catalyst is used with a large amount, a solvent may not be specifically added. In this case, an amount thereof is 5 to 100-fold amount, preferably 10 to 30-fold amount by weight based on the amount of Compound (XXVI).

The reaction is carried out in the temperature range
25 usually at 0 to 250°C, preferably at 100 to 200°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 10 minutes to 12 hours, preferably 30 minutes to 5 hours.

A step of obtaining Compound (XXVIII) by brominating
30 Compound (XXVII) in Step H4 can be easily carried out by the conventionally known method, for example, in which reaction is carried out under irradiation of light using N-bromosuccineimide as a brominating agent, or reaction is carried out in the presence of a radical initiating agent
35 such as azobisisobutyronitrile (AIBN), benzoyl peroxide (BPO), etc. The solvent to be used is not specifically

limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, halogenated hydrocarbons such as dichloromethane, chloroform, carbon tetrachloride, 1,2-dichloroethane, etc., preferably 1,2-dichloroethane.

When the reaction is carried out under irradiation of light, as a light source, a mercury lamp is suitably used. The N-bromosuccinimide to be used is usually 1 to 2.5-fold mol amount, preferably 1 to 1.1-fold mol amount based on the amount of Compound (XXVII). A reaction temperature is usually in the range of 0 to 80°C, preferably 30 to 60°C.

A reaction time may vary depending on the reaction temperature and the like, and it is usually 1 hour to 12 hours, preferably 3 hours to 8 hours.

When the reaction is carried out in the presence of a radical initiator such as azobisisobutyronitrile (AIBN), benzoyl peroxide (BPO), etc., an amount of the radical initiator to be used is usually 0.001 to 0.1-fold mol amount, preferably 0.01 to 0.05-fold mol amount based on the amount of Compound (XXVII).

The N-bromosuccinimide to be used is usually 1 to 4-fold mol amount, preferably 1 to 1.1-fold mol amount based on the amount of Compound (XXVII).

The reaction is carried out in the temperature range usually at 50 to 100°C, preferably at 60 to 80°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 30 minutes to 12 hours, preferably 1 hour to 6 hours.

A step of obtaining Compound (XXIX) by removing hydrogen bromide from Compound (XXVIII) in Step H5 is carried out by treating Compound (XXVIII) with a base (preferably triethylamine or 1,5-diazabicyclo[4.3.0]-5-nonene (DBN)).

The solvent to be used is not specifically limited

so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, halogenated hydrocarbons such as dichloromethane, carbon tetrachloride, 1,2-dichloroethane, etc.; aromatic hydrocarbons such as benzene, toluene, xylene, etc.; esters such as methyl acetate, ethyl acetate, etc.; nitriles such as acetonitrile, propionitrile, etc., preferably 1,2-dichloroethane, toluene, ethyl acetate or acetonitrile.

10 An amount of the base to be used is 1 to 150-fold mol amount, preferably 1 to 15-fold mol amount based on the amount of Compound (XXVIII).

 The reaction is carried out in the temperature range usually at 0 to 150°C, preferably at 60 to 120°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 30 minutes to 12 hours, preferably 1 hour to 3 hours.

 Preparation process I is another process for producing Compound (XXVI) in Preparation process H.

20 Bromination reaction of Compound (XXX) in Step I1 is carried out in the same manner as in Step H4 except for using Compound (XXX) in place of Compound (XXVII).

 Step I2 can be easily carried out by reacting Compound (XXXI) and triphenylphosphine in a solvent.

25 The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, halogenated hydrocarbons such as dichloromethane, chloroform, carbon tetrachloride, 1,2-dichloroethane, etc.; aromatic hydrocarbons such as benzene, toluene, xylene, etc.; esters such as methyl acetate, ethyl acetate, etc.; nitriles such as acetonitrile, propionitrile, etc.; amides such as N,N-dimethylformamide, N,N-dimethylacetamide, N-methylpyrrolidone, etc., ketones such as acetone, methyl ethyl ketone, etc.; or dimethylsulfoxide, preferably ethyl acetate, or aceto-

30

35

nitrile.

An amount of the triphenylphosphine to be used is 1 to 2-fold mol amount, preferably 1 to 1.5-fold mol amount based on the amount of Compound (XXXI).

5 The reaction is carried out in the temperature range usually at 0 to 150°C, preferably at 60 to 100°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 30 minutes to 12 hours, preferably 1 hour to 3 hours.

10 A reaction of Compound (XXXII) and Compound (XXXIII) in Step I3 is a reaction known as the so-called Wittig reaction, and carried out optionally by selecting the conventionally known conditions.

15 An amount of Compound (XXXIII) to be used is usually 1 to 10-fold mol amount, preferably 1 to 1.5-fold mol amount based on the amount of Compound (XXXII).

20 The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, halogenated hydrocarbons such as dichloromethane, carbon tetrachloride, chloroform, 1,2-dichloroethane, etc.; ethers such as diethyl ether, tetrahydrofuran, dioxane, dimethoxyethane, etc.; aromatic hydrocarbons such as benzene, toluene, xylene, etc.; esters
25 such as methyl acetate, ethyl acetate, etc.; nitriles such as acetonitrile, propionitrile, etc.; amides such as N,N-dimethylformamide, N,N-dimethylacetamide, N-methylpyrrolidone, etc.; dimethylsulfoxide or a mixed solvent of the above-mentioned solvents, and further, a mixed solvent of
30 the above solvents and water may be also used.

35 The base to be used may include alkali metal hydroxides such as lithium hydroxide, sodium hydroxide, potassium hydroxide, etc.; alkali metal carbonates such as sodium carbonate, potassium carbonate, etc.; alkali metal alkoxides such as sodium methoxide, sodium ethoxide, sodium t-butoxide, potassium t-butoxide, etc.; alkali metal

hydrides such as sodium hydride, potassium hydride, etc.; alkyl lithium such as methyl lithium, butyl lithium, etc.; metal amides such as sodium amide, lithium diisopropyl amide, etc.; and organic amines such as triethylamine, diisopropylethylamine, tripropylamine, 1,5-diazabicyclo-
5 [4.3.0]-5-nonene (DBN), etc. An amount of the base to be used is usually 1 to 5-fold mol amount, preferably 1 to 2-fold mol amount based on the amount of is Compound (XXXII).

The reaction is carried out in the temperature range
10 usually at 0 to 150°C, preferably at 15 to 80°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 1 hour to 24 hours, preferably 1 hour to 6 hours.

Incidentally, in this step, a mixture of cis and
15 trans geometric isomer can be usually obtained and it can be applied to the next step in the state of a mixture.

Hydrolysis reaction of Compound (XXXIV) to the carboxyl group in Step I4 may be also carried out, for example, by a typical alkali hydrolysis reaction using
20 sodium hydroxide or potassium hydroxide, or a method disclosed in "Protective Group in Organic Synthesis" written by W. Greene and P.G.H. Wult, 2nd Ed., John Wiley & Sons, p. 229.

Reduction of Compound (XXXV) in Step I5 can be
25 carried out by a catalytic hydrogenation reaction or a reduction using hydroiodic acid/red phosphorus, and when W is a halogen atom, the reduction using hydroiodic acid/red phosphorus is suitable.

In the case of the catalytic hydrogenation reaction,
30 Compound (XXXV) is subjected to catalytic reduction by hydrogen in the presence of a catalyst whereby Compound (XXVI) can be synthesized.

The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dis-
35 solve the starting materials with a certain extent, and there may be mentioned, for example, alcohols such as

methanol, ethanol, etc., or ethers such as tetrahydrofuran, dioxane, etc., preferably alcohols.

As the catalyst to be used, there may be mentioned, for example, palladium-carbon, platinum-carbon, platinum
5 black, rhodium-carbon or Raney nickel, preferably palladium-carbon.

In the catalytic hydrogenation reaction, a partial pressure of hydrogen is usually 1 atm. to 10 atm., preferably 1 atm. to 5 atm.

10 The reaction is carried out in the temperature range usually at 0°C to 100°C, preferably at 20°C to 80°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 15 minutes to 72 hours, preferably 30 minutes to 48 hours.

15 In the case of the reduction using hydroiodic acid/red phosphorus, the solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, water, acetic acid or
20 propionic acid, preferably acetic acid. The hydroiodic acid to be used is 3 to 15-fold mol amount, preferably 4 to 8-fold mol amount based on 1 mol of Compound (XXXV).

The red phosphorus to be used is 1 to 10-fold mol amount, preferably 2 to 5-fold mol amount based on the
25 amount of Compound (XXXV). A method of addition thereof may be once or divided into several times.

The reaction is carried out in the temperature range usually at 90 to 150°C, preferably at 100 to 130°C. A reaction time may vary depending on the reaction tempera-
30 ture and the like, and it is usually 1 hour to 48 hours, preferably 6 hours to 24 hours.

Preparation process J is a process for producing Compound (XVI) by demethylating Compound (XXXVI). Demethylation of Compound (XXXVI) is carried out by
35 optionally employing a conventionally known method (see "Protective Group in Organic Synthesis" written by W.Greene

and P.G.H. Wult, 2nd Ed., John Wiley & Sons, p. 146), for example, a method of using hydrobromic acid or hydroiodic acid, or a method of using boron tribromide and trimethylsilane iodide, and the like.

5 Preparation process K is a preparation process of Compound (XVIII), and carried out easily by employing a conventionally known method, for example, a method as disclosed in J.Org.Chem., 61, 3398 (1996).

10 Preparation process M is another process for producing Compound (XXVI).

 Step M1 is a process for producing Compound (XXXV') by condensation reaction of Compound (XXX) and Compound (XXXIII) in the presence of a base.

15 The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, alcohols such as methanol, ethanol, propanol, butanol, t-butanol, etc.; ethers such as tetrahydrofuran, dioxane, 1,2-dimethoxyethane, etc.; amides such as N,N-dimethylformamide, N,N-dimethylacetamide, N-methylpyrrolidone, etc., preferably
20 N,N-dimethylformamide.

 An amount of Compound (XXXIII) to be used is usually 1 to 1.5-fold mol amount, preferably 1 to 1.2-fold mol
25 amount based on the amount of Compound (XXX).

 The base to be used may be mentioned alkali metal alkoxides such as sodium methoxide, sodium ethoxide, sodium t-butoxide, potassium t-butoxide, etc.; or alkali metal hydrides such as sodium hydride, potassium hydride, etc.,
30 preferably sodium t-butoxide. An amount thereof to be used is usually 1 to 2-fold mol amount, preferably 1 to 1.5-fold mol amount based on the amount of Compound (XXX).

 The reaction is carried out in the temperature range usually at 0 to 100°C, preferably at 10 to 40°C. A reaction time may vary depending on the reaction temperature
35 and the like, and it is usually 1 to 24 hours, preferably 3

to 12 hours.

Step M2 is a hydrogenation reaction of olefin and can be carried out in the same manner as in Step I5 of Preparation process I.

5 Preparation process N is another process for producing Compound (XXIX).

Step N1 is a step of obtaining Compound (XXXX) by reacting Compound (XXVII) and N-bromosuccinimide under irradiation of light, preferably under irradiation of light
10 using a mercury lamp light source.

The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, halogenated hydrocarbons such as dichloromethane, chloroform, carbon tetra-
15 chloride, 1,2-dichloroethane, etc., preferably 1,2-dichloroethane. An amount of the N-bromosuccinimide to be used is usually 1.8 to 2.5-fold mol amount, preferably 2.0 to 2.1-fold mol amount based on the amount of Compound
20 (XXVII).

The reaction is carried out in the temperature range usually at 0 to 80°C, preferably at 30 to 60°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 1 hour to 12 hours, preferably 3
25 hours to 8 hours.

Step N2 is a step of converting from vicinal dibromide to olefin and is carried out easily by a method as disclosed in, for example, Synth., Comm., 26, 3791 (1996) or the like.

30 Preparation process O is a preparation process of an intermediate Compound (XXXXX) for the exemplary Compounds (If) and (Ig) wherein X=a methylene group among the compounds represented by the formula (I).

Step O1 is a step of producing Compound (XXXXIII) by
35 reacting an acetylene compound (XXXXII) to Compound (XXXXI).

The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be preferably mentioned, for example, ethers such as tetrahydrofuran, dioxane, 1,2-dimethoxyethane, etc.

An amount of the acetylene compound (XXXXII) to be used is usually 1 to 2-fold mol amount, preferably 1 to 1.5-fold mol amount based on the amount of Compound (XXXXI).

The reaction is carried out in the temperature range usually at 0 to 100°C, preferably at 0 to 30°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 5 minutes to 12 hours, preferably 30 minutes to 6 hours.

The reaction of producing Compound (XXXXIV) from Compound (XXXXIII) in Step 02 is carried out by subjecting to conversion reaction in a solvent in the presence of an acid catalyst.

The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and there may be mentioned, for example, alcohols such as methanol, ethanol, propanol, butanol, etc.; ethers such as tetrahydrofuran, dioxane, 1,2-dimethoxyethane, etc.; nitriles such as acetonitrile, etc.; and a mixed solvent of amides such as N,N-dimethylformamide, N,N-dimethylacetamide, N-methylpyrrolidone, etc. and water, preferably methanol, ethanol, tetrahydrofuran or a mixed solvent of N,N-dimethylformamide and water.

As the acid catalyst to be used, there may be mentioned, for example, mineral acids such as hydrochloric acid, sulfuric acid, phosphoric acid, etc.; or organic acids such as methanesulfonic acid, trifluoroacetic acid, etc., preferably trifluoroacetic acid. An amount of the catalyst to be used is usually 1 to 100-fold mol amount, preferably 1 to 50-fold mol amount based on the amount of

Compound (XXXXIII).

The reaction is carried out in the temperature range usually at 0 to 100°C, preferably at 0 to 30°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 5 minutes to 48 hours, preferably 30 minutes to 24 hours.

Step 03 is a step of obtaining Compound (XXXXVI) by reacting Compound (XXXXIV) and Compound (XXXXV) in the presence of a base in a solvent.

10 An amount of Compound (XXXXV) to be used is usually 1 to 10-fold mol amount, preferably 1 to 5-fold mol amount based on the amount of Compound (XXXXIV).

As the base to be used, there may be mentioned, for example, alkali metal hydrides such as sodium hydride, 15 lithium hydride, etc.; alkali metal alkoxides such as sodium methoxide, sodium ethoxide, potassium t-butoxide, etc.; alkyl lithiums such as methyl lithium, butyl lithium, etc.; metal amides such as sodium amide, lithium diisopropyl amide, etc., preferably alkali metal hydride. An 20 amount of the base to be used is usually 1 to 5-fold mol amount, preferably 1 to 2-fold mol amount based on the amount of Compound (XXXXV).

The solvent to be used is not specifically limited so long as it does not inhibit the reaction and can dissolve the starting materials with a certain extent, and 25 there may be mentioned, for example, aromatic hydrocarbons such as benzene, toluene, etc.; ethers such as tetrahydrofuran, dioxane, dimethoxyethane, dimethoxyethane, etc.; or aprotic polar solvents such as N,N-dimethylformamide, N,N-dimethylacetamide, N-methylpyrrolidone, dimethylsulfoxide, 30 etc., preferably ethers.

The reaction is carried out in the temperature range usually at -50°C to 100°C, preferably at -10°C to 50°C. A reaction time may vary depending on the reaction temperature and the like, and it is usually 15 minutes to 12 35 hours, preferably 30 minutes to 5 hours.

In Step 04, Compound (XXXXVII) can be synthesized by
subjecting Compound (XXXXVI) to catalytic reduction with
hydrogen in a solvent in the presence of a base, and the
step is carried out in the same manner as in the catalytic
5 hydrogenation of Step I5.

Step 05 is a conversion to olefin by dehydrogenation
and, for example, it is carried out in the same manner as
in the above-mentioned Preparation process N except for
using Compound (XXXXVII) in place of Compound (XXVII).

10 Step 06 is demethylation of a methoxy group and
hydrolysis of an ester group.

Demethylation of the methyl group and hydrolysis of
the ester group are simultaneously accomplished by carrying
out in the same manner as in, for example, Preparation
15 process J except for using Compound (XXXXVIII) in place of
Compound (XXXVI).

Step 07 is a protection step for a carboxyl group by
esterification, and can be easily converted by a conven-
tionally known method (for example, see "Protective Group
20 in Organic Synthesis" written by W. Greene and P.G.H. Wult,
2nd Ed., John Wiley & Sons, p. 224).

Preparation process P is another process for produc-
ing Compound (XVIII).

Step P1 is a step of producing Compound (XXXXXI) by
25 reacting Compound (XXI) and triphenylphosphine, and Step P2
is a step of producing Compound (XVIII) by reacting Com-
pound (XXXXXI) and formaldehyde in the presence of a base.

Step P1 is carried out in the same manner as in Step
I2 except for using Compound (XXI) in place of Compound
30 (XXI).

Step P2 is carried out in the same manner as in Step
I3 except for using Compound (XXXXXI) in place of Compound
(XXXII), and using formaldehyde in place of Compound
(XXXIII).

35 The compound represented by the present invention
has a potent leukotriene antagonistic action, and extremely

useful as an antiazma agent, an antiallergic agent and anti-inflammatory agent.

When used as a prophylactic or therapeutic medication for the diseases described above, Compound (I) or a
5 pharmaceutically acceptable salt thereof of the present invention can be administered alone or can be presented as part of a pharmaceutical formulation. The pharmaceutical formulation is prepared by blending the active ingredient with appropriate pharmaceutically acceptable excipient,
10 diluent, etc., followed by formulation in the form of tablets, capsules, granules, powders or syrups, etc. for oral administration or in the form of injections, etc. for parenteral administration (preferably oral administration).

The production of such pharmaceutical formulation is
15 carried out according to general technique known to those skilled in the art using additives such as an excipient (for example, a sugar derivative such as lactose, sucrose, glucose, mannitol, or sorbitol; a starch derivative such as corn starch, potato starch, α -starch, dextrin, or carboxy-
20 methyl starch; a cellulose derivative such as crystalline cellulose, low-substituted hydroxypropyl cellulose, hydroxypropylmethyl cellulose, carboxymethyl cellulose, calcium carboxymethyl cellulose, or internally bridged sodium carboxymethyl cellulose; acacia; dextran; pullulan;
25 a silicate derivative such as light silicic acid anhydride, synthetic aluminum silicate, or magnesium aluminate meta-silicate; a phosphate derivative such as calcium phosphate; a carbonate derivative such as calcium carbonate; a sulfate derivative such as calcium sulfate; etc.), a binder (for
30 example, one of the excipients described above; gelatin; polyvinylpyrrolidone; Macrogol (tradename), etc.), a disintegrator (for example, one of the excipients described above; a chemically modified starch, cellulose derivative such as sodium croscarmellose, sodium carboxymethyl starch;
35 bridged polyvinylpyrrolidone, etc.), a lubricant (for example, talc; stearic acid; a metal salt of stearic acid

such as calcium stearate, or magnesium stearate; colloidal silica; a wax such as bee gum and spermaceti; boric acid; glycol; a carboxylic acid such as fumaric acid, or adipic acid; a sodium carboxylate such as sodium benzoate; a sulfate such as sodium sulfate; leucine; a laurylsulfate such as sodium laurylsulfate, or magnesium laurylsulfate; a silicic acid such as silicic acid anhydride, or a silicic acid hydrate; one of the starch derivatives described above in relation to the excipient, etc.), a stabilizer (for example, a p-hydroxybenzoate derivative such as methylparaben, or propylparaben; an alcohol such as chlorobutanol, benzyl alcohol, or phenylethyl alcohol; benzalkonium chloride; a phenol derivative such as phenol, or cresol; thimerosal; acetic anhydride; sorbic acid, etc.), a corrigent (for example, a sweetening, souring, or flavoring agent, which are conventionally used, etc.), a diluent, a solvent for injection (for example, water, ethanol, glycerin, etc.), etc. A dose for administration may vary depending on symptom, an age, etc, and in the case of oral administration, it is desirably administered 1 mg (preferably 5 mg) as a lower limit and 1000 mg (preferably 500 mg) as an upper limit once per day, and in the case of intravenous administration, it is desirably administered 0.1 mg (preferably 1mg) as a lower limit and 500 mg (preferably 300 mg) as an upper limit once per day to an adult person with one to six times per day depending on the symptom.

Utilizability in industry

The compound represented by the formula (I) or a pharmaceutically acceptable salt thereof according to the present invention has leukotriene C₄ antagonistic action and leukotriene E₄ antagonistic action in addition to potent leukotriene D₄ antagonistic action, and is extremely useful as an antiasthmatic agent, an antiallergic agent and an anti-inflammatory agent.

Example

In the following, the present invention is explained by referring to Test examples and Examples, but the scope of the present invention is not limited to these.

5 Test example 1

Leukotriene D₄ receptor binding test

<Preparation of receptor sample>

As a receptor sample, a lung cell membrane fraction from guinea pigs was used. Preparation of the membrane
10 fraction was carried out according to the method of Ahn et al. (Eur. J. Pharmacol., 127, 153-155 (1986)). Lungs of Hartley male guinea pigs (400 to 500 g body weight, Japan SLC Inc.) were isolated, and perfused with a physiological saline, and then, adding 10 mM of PIPES, 10 mM of MgCl₂ and
15 10 mM of CaCl₂ buffer (pH 7.5) to the lung tissue and the mixture was homogenized. This homogenate was centrifuged at 70,000xg for 10 minutes to obtain a membrane fraction.
<Leukotriene D₄ receptor binding test>

Leukotriene D₄ (LTD₄) receptor binding test was
20 carried out according to the method of Aharony, et al. (J. Pharmacol. Expl. Ther., 243, 921-926 (1987)). To 0.42 mg of the receptor sample were added 10 mM of PIPES, 10 mM of MgCl₂ and 10 mM of CaCl₂ buffer (pH 7.5) to make the total amount of 480 μ l, and 10 μ l of [³H] LTD₄ (NEN Life Science
25 Products Inc.) and 10 μ l of a test compound in dimethylsulfoxide were added to the mixture, and the resulting mixture was incubated at 25°C for 30 minutes. The mixtures thus incubated were filtered through a glass fiber filter (Whatman International Ltd., GF/C) using cell harvester
30 (Biomedical Research & Development Laboratories, Inc., M-30R). The filter paper was washed with 10 mM of Tris and 100 mM of NaCl buffer (pH 7.5), and 5 ml of a liquid scintillator (nacalai tesque inc., clearsol I) was added thereto, and radioactivity was measured by a liquid
35 scintillation analyzer (Packard Instrument Co., 2000CA). When a dissociation constant (K_d) of LTD₄ was to be

obtained, [³H] LTD₄ with 0.03 to 0.5 nM was used, and 1 μM of non-radioactive LTD₄ was added. When a binding inhibition constant (K_i) of the Test compound is to be measured, [³H] LTD₄ with 0.2 nM was used. K_d and K_i are calculated according to the method of Bennett et al. (Neurotransmitter Receptor Binding, 2nd ed., edited by H.I. Yamamura et al., pp. 61-89, Raven Press (1985)).

Table 2
Results of leukotriene D₄ receptor binding test

Compound to be tested	pK _i
Example Compound 3	10.0
Example Compound 4	9.7
Example Compound 7	9.7
Example Compound 9	9.8
Example Compound 10	10.0
Example Compound 11	9.9
Example Compound 12	9.7
Example Compound 13	9.9
Example Compound 17	9.8
Example Compound 29	9.6
Example Compound 30	9.7
Example Compound 31	9.7
Compound A	9.5

Compound A; 3-[2-(7-chloro-6-fluoroquinolin-2-yl)methoxy-6,11-dihydrobenz[b,e]oxepin-11-yl]thiopropionic acid (see WO 94/193445 publication)

Test example 2

Leukotriene D₄ induced respiratory constriction test
Respiratory constriction was measured by modifying the method of Konzett and Rössler (Arch. Exp. Pathol. Pharmacol., 195, pp. 71-74 (1940)). Hartley male guinea

pigs (400 to 500 g body weight, Japan SLC Inc.) were anesthetized with pentobarbital (50 mg/kg, s.c.), and a cannula was inserted into the trachea to carry out artificial ventilation with an artificial ventilator (manufactured by Harvard Apparatus, Model 683). An inner pressure of the respiratory tract was measured by a differential pressure transducer (Nihon Kohden, TP-603T) connected to the respiratory cannula and it is used as an index of respiratory constriction.

10 LTD₄ (0.03, 0.06, 0.13, 0.25, 0.5, 1 and 2 µg/kg, Sigma-Aldrich) was intravenously administered from a cannula inserted into the right jugular vein from a low dose with an interval of 5 minutes to cause a respiratory constricttion reaction and an increased amount of a
15 respiratory inner pressure was measured. Test compound was suspended in 0.5% sodium carboxymethyl cellulose aqueous solution or dissolved in 30% propylene glycol aqueous solution, and orally administered 24 hours before administration of LTD₄. Animals were fasted for 24 hours
20 before administration of the Test compound. From a dose-reaction curve of LTD₄, 50% reaction dose (ED₅₀) was measured, and a dose of the Test compound required for shifting two-times of a dose-reaction curve of a control group to a higher dose side was calculated from the formula
25 shown below and it was made as LTD₄ inhibitory activity.
LTD₄ inhibitory activity = (Dose of Compound administered) / { (ED₅₀ of group to which Compound was added) / (ED₅₀ of control group) - 1 }

Table 3
Results of leukotriene D₄ induced
respiratory constriction test

Compound to be tested	LTD ₄ inhibitory activity (µg/kg, p.o. 24h)
Example Compound 3	4.71
Example Compound 4	1.91
Example Compound 7	0.85
Example Compound 8	5.30
Example Compound 10	9.93
Example Compound 14	4.85
Example Compound 16	0.79
Example Compound 18	1.12
Example Compound 19	2.43
Example Compound 20	0.70
Example Compound 28	1.70
Compound A	25.52

5

Compound A; 3-[2-(7-chloro-6-fluoroquinolin-2-yl)-methoxy-6,11-dihydrobenz[b,e]oxepin-11-yl]thiopropionic acid (see WO 94/193445 publication)

10 Preparation example 1

Tablet

Compound of Example 7	30.0 mg
Lactose	144.0
Corn starch	25.0
Magnesium stearate	1.0
	200 mg

15

A tablet is prepared using the ingredients above. The components are blended and compressed by a tablet machine to form a tablet weighing 200 mg. The tablet may be coated if necessary, for example, to form a sugar-coated tablet or a coated tablet.

20

Example 1

Methyl [3-[(E)-2-(6,7-difluoroquinolin-2-yl)-ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetate (methyl ester of Exemplary compound 1)

5 After a solution of 1.19 g (2.99 mmol) of 3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-ol dissolved in 10 ml of tetrahydrofuran was cooled to 0°C with ice, 0.85 ml (5.98 mmol) of triethylamine and 0.30 ml (3.89 mmol) of methanesulfonyl chloride
10 were added to the solution, and the mixture was stirred at 0°C for 1 hour, and further at room temperature for 3 hours.

 After completion of the reaction, the solvent was removed from the mixture under reduced pressure. The
15 residue was dissolved in 15 ml of N,N-dimethylformamide, then, 0.54 g (5.98 mmol) of methyl glycolate was added to the mixture, and the resulting mixture was stirred at room temperature overnight.

 After completion of the reaction, water was added to
20 the reaction mixture, and extracted with chloroform. The organic layer was washed with water, dried over anhydrous sodium sulfate, and the solvent was removed under reduced pressure. The residue was applied to silica gel column chromatography (solvent: hexane/ethyl acetate=2/1 (volume
25 ratio)) to obtain 0.38 g of the title compound as yellowish solid.

EI-MS (m/z); 469 (M^+).

CI-MS (m/z); 470 (M^++1).

Example 2

30 [3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid (Exemplary compound 1)

 To mixed solution comprising 15 ml of methanol and 5 ml of tetrahydrofuran was dissolved 0.38 g (0.81 mmol) of
35 methyl [3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetate, 2.4 ml (2.4 mmol)

of an aqueous 1N sodium hydroxide solution was added to the solution and the resulting mixture was stirred at room temperature for 5 hours.

After completion of the reaction, the reaction mixture was adjusted to pH 6.5 by using an aqueous diluted acetic acid solution, and then, the mixture was concentrated under reduced pressure. Water was added to the residue, and the precipitated solid was collected by filtration. The solid was applied to silica gel column chromatography (solvent: chloroform/methanol=9/1 (volume ratio)) to obtain 0.21 g of the title compound as yellowish solid.

FAB-MS (m/z); 456 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 4.14 (s, 2H), 4.93 (s, 1H), 7.18-7.67 (m, 9H), 7.89-8.07 (m, 5H), 8.37 (d, $J=8.5\text{Hz}$, 1H).

Example 3

Sodium [3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetate (sodium salt of Exemplary compound 1)

In a mixed solution of 15 ml of tetrahydrofuran and 5 ml of methanol was dissolved 0.20 g (0.44 mmol) of [3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid, 0.44 ml (0.44 mmol) of an aqueous 0.1N sodium hydroxide solution was added to the solution and the mixture was stirred at room temperature for 1 hour.

After completion of the reaction, the reaction mixture was concentrated, ethanol was added to the residue, and the precipitated solid was collected by filtration to obtain 0.14 g of the title compound as pale yellowish solid.

m.p.; 213 to 227°C.

FAB-MS (m/z); 478 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 3.81 (s, 2H), 5.03 (s, 1H), 7.15-7.65 (m, 9H), 7.89-8.06 (m, 5H), 8.37 (d, $J=8.8\text{Hz}$, 1H).

Example 4

Sodium 3-{[3-[(E)-2-(6,7-difluoroquinolin-2-yl)-ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionate (sodium salt of Exemplary compound 8)

(a) In a mixed solution of 0.4 ml of trifluoroacetic acid and 40 ml of methylene chloride was dissolved 0.33 g (0.83 mmol) of 3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-ol, 0.08 ml (0.92 mmol) of 3-mercaptopropionic acid was added to the solution and the mixture was stirred at room temperature for 3 hours.

After completion of the reaction, the reaction mixture was concentrated, and the residue was applied to silica gel column chromatography (eluent: chloroform/methanol=19/1 (volume ratio)) to obtain 0.17 g (0.35 mmol) of 3-{[3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid as brown solid.

(b) 3-{[3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid obtained in Example 4(a) was dissolved as such in a mixed solution of 10 ml of tetrahydrofuran and 20 ml of methanol, 0.35 ml (0.35 mmol) of an aqueous 1N sodium hydroxide solution was added to the solution and the mixture was stirred at room temperature for 1 hour.

After completion of the reaction, the reaction mixture was concentrated, the residue was washed with diethyl ether, and dried under reduced pressure to obtain 0.14 g of the title compound as pale yellowish solid. m.p.; 213 to 216°C.

FAB-MS (m/z); 508 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 1.90-2.10 (m, 2H), 2.30-2.45 (m, 2H), 5.48 (s, 1H), 7.00 (s, 2H), 7.20-7.60 (m, 6H), 7.33 (d, $J=8.8\text{Hz}$, 1H), 7.37 (d, $J=7.3\text{Hz}$, 1H), 7.48 (d, $J=8.5\text{Hz}$, 1H), 7.54 (d, $J=16.4\text{Hz}$, 1H), 7.66 (d, $J=7.8\text{Hz}$, 1H), 7.80-8.20 (m, 3H), 7.87 (d, $J=16.1\text{Hz}$, 1H), 8.04 (dd, $J=9.0\text{Hz}$, 2.2Hz, 1H), 8.38 (d, $J=8.8\text{Hz}$, 1H).

Reaction was carried out in the same manner as in

Example 4 to obtain compounds of the following Examples 5 to 9.

Example 5

Sodium 3-([3-(7-chloro-6-fluoroquinolin-2-yl)-
5 methoxy-5H-dibenzo[a,d]cyclohepten-5-yl]thio)propionate
(sodium salt of Exemplary compound 718)

Appearance; pale yellowish solid

m.p.; 229 to 232°C.

¹H-NMR (δ, DMSO-d₆); 2.00-2.20 (m, 2H), 2.25-2.45 (m, 2H),
10 5.38 (s, 1H), 5.41 (s, 2H), 6.83 (d, J=12.0Hz, 1H), 6.90
(d, J=12.2Hz, 1H), 7.00 (dd, J=8.5Hz, 2.7Hz, 1H), 7.20-7.55
(m, 6H), 7.75 (d, J=8.5Hz, 1H), 8.05 (d, J=9.8Hz, 1H), 8.29
(d, J=7.3Hz, 1H), 8.44 (d, J=8.3Hz, 1H).

Example 6

15 Sodium 3-([3-[(E)-2-(6,7-difluoroquinolin-2-yl)-
ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio)-2-(S)-
methylpropionate (sodium salt of Exemplary compound 9)

Appearance; pale yellowish solid

m.p.; 240 to 249°C.

20 FAB-MS (m/z); 522 (M⁺+1).

¹H-NMR (δ, DMSO-d₆); 0.89 (d, J=6.6Hz, 3H), 2.10-2.89 (m,
2H), 2.51-2.58 (m, 1H), 5.45 (s, 1H), 7.00 (s, 2H), 7.31-
7.49 (m, 5H), 7.54 (d, J=16.4Hz, 1H), 7.66 (d, J=8.1Hz,
1H), 7.85 (d, J=6.8Hz, 1H), 7.86-7.98 (m, 2H), 7.98 (d,
25 J=3.9Hz, 1H), 8.04 (dd, J=11.0Hz, 8.8Hz, 1H), 8.38 (d,
J=8.8Hz, 1H).

Example 7

Sodium 3-([3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio)propionate
30 (sodium salt of Exemplary compound 478)

Appearance; yellowish solid

m.p.; 228 to 238°C.

FAB-MS (m/z); 524 (M⁺+1).

¹H-NMR (δ, DMSO-d₆); 1.96-2.01 (m, 2H), 2.34-2.37 (m, 2H),
35 5.48 (s, 1H), 7.00 (s, 2H), 7.29-7.50 (m, 5H), 7.55 (d,
J=16.4Hz, 1H), 7.66 (dd, J=8.8Hz, 1.5Hz, 1H), 7.88 (d,

J=16.4Hz, 1H), 7.89 (d, J=1.5Hz, 1H), 8.00 (d, J=8.8Hz, 1H), 8.01 (d, J=10.0Hz, 1H), 8.23 (d, J=7.1Hz, 1H), 8.39 (d, J=8.5Hz, 1H).

Example 8

5 Sodium 3-{{[3-[(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionate (sodium salt of Exemplary compound 1418)

Appearance; orange solid

m.p.; 193 to 203°C.

10 FAB-MS (m/z); 506 (M⁺+1).

¹H-NMR (δ, DMSO-d₆); 2.02-2.08 (m, 2H), 2.35-2.41 (m, 2H), 5.49 (s, 1H), 7.00 (s, 2H), 7.29-7.49 (m, 5H), 7.53-7.61 (m, 2H), 7.67 (d, J=8.1Hz, 1H), 7.86-7.94 (m, 3H), 7.99-8.03 (m, 2H), 8.41 (d, J=8.5Hz, 1H).

15 Example 9

Sodium 3-{{[3-[(E)-2-(5,6,7,8-tetrahydroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionate (sodium salt of Exemplary compound 1885)

Appearance; pale yellowish solid

20 m.p.; 175 to 185°C.

FAB-MS (m/z); 508 (M⁺+1).

¹H-NMR (δ, DMSO-d₆); 1.76-1.85 (m, 4H), 2.08-2.13 (m, 2H), 2.35-2.40 (m, 2H), 2.74-2.84 (m, 4H), 5.44 (s, 1H), 6.97 (s, 2H), 7.28-7.46 (m, 8H), 7.54-7.59 (m, 2H), 7.78 (s, 25 1H).

Reaction was carried out in the same manner as in Examples 1 to 3 to obtain compounds of the following Examples 10 to 13.

Example 10

30 Sodium 3-{{[3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}-2-(R)-methylpropionate (sodium salt of Exemplary compound 4)

Appearance; yellowish solid

¹H-NMR (δ, DMSO-d₆); 0.74, 1.20 (each b, 3H in total), 2.44 (b, 1H), 3.53, 3.77 (each b, 2H in total), 4.68, 5.55 (each s, 1H in total), 7.00-7.70 (m, 9H), 7.87-8.07 (m, 5H), 8.37

(d, J=8.8Hz, 1H).

Example 11

Sodium [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetate (sodium salt of Exemplary compound 471)

Appearance; yellowish solid

¹H-NMR (δ, DMSO-d₆); 3.81 (bs, 2H), 5.04 (bs, 1H), 7.15 (bs, 2H), 7.26-7.66 (m, 7H), 7.90-8.02 (m, 4H), 8.22 (d, J=7.3Hz, 1H), 8.38 (d, J=8.6Hz, 1H).

Example 12

Sodium 3-{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}propionate (sodium salt of Exemplary compound 473)

Appearance; yellowish solid

¹H-NMR (δ, DMSO-d₆); 2.37 (bs, 2H), 3.72, 3.82 (each bs, 2H in total), 4.70 (bs, 1H), 7.18 (bs, 2H), 7.28-7.66 (m, 8H), 7.88-8.04 (m, 4H), 8.27 (d, J=7.1Hz, 1H), 8.38 (d, J=8.5Hz, 1H).

Example 13

Sodium 3-{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}-2-(R)-methylpropionate (sodium salt of Exemplary compound 474)

Appearance; yellowish solid

¹H-NMR (δ, DMSO-d₆); 0.73, 1.20, 1.52 (each b, 3H in total), 2.45 (b, 1H), 3.33-3.86 (each m, 2H), 4.68, 5.55 (each s, 1H in total), 7.19-7.63 (m, 9H), 7.88-8.27 (m, 5H), 8.38 (d, J=8.8Hz, 1H).

Reaction was carried out in the same manner as in Example 4 to obtain compounds of the following Examples 14 to 16.

Example 14

Sodium [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thioacetate (sodium salt of Exemplary compound 476)

Appearance; orange solid

¹H-NMR (δ, DMSO-d₆); 2.61 (d, J=13.7Hz, 2H), 2.66 (d,

J=13.7Hz, 2H), 5.63 (s, 1H), 6.98 (s, 2H), 7.31 (td, J=9.0Hz, 2.0Hz, 1H), 7.33-7.42 (m, 3H), 7.45 (d, J=7.8Hz, 1H), 7.52 (d, J=16.4Hz, 1H), 7.67 (dd, J=8.1Hz, 1.5Hz, 1H), 7.75 (d, J=2.0Hz, 1H), 7.88 (d, J=16.1Hz, 1H), 7.98 (d, J=9.8Hz, 1H), 8.02 (d, J=10.0Hz, 1H), 8.24 (d, J=7.6Hz, 1H), 8.39 (d, J=8.6Hz, 1H).

Example 15

Sodium 2-{{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionate
 10 (sodium salt of Exemplary compound 477)

Appearance; pale yellowish solid

¹H-NMR (δ, DMSO-d₆); 1.10-1.14 (m, 3H), 2.68-2.76 (m, 3H), 5.67 (d, J=6.4Hz, 1H), 6.98 (s, 2H), 7.23-7.54 (m, 6H), 7.67 (dd, J=8.1Hz, 1.5Hz, 1H), 7.31-7.90 (m, 2H), 7.98 (d, J=7.6Hz, 1H), 8.01 (d, J=9.8Hz, 1H), 8.23 (d, J=7.3Hz, 1H), 8.39 (d, J=8.6Hz, 1H).

Example 16

Sodium 3-{{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thiomethyl}-cyclopropaneacetate (sodium salt of Exemplary compound 487)
 20

Appearance; pale yellowish solid

¹H-NMR (δ, DMSO-d₆); 0.15 (bs, 2H), 0.33 (bs, 2H), 1.95 (d, J=14.2Hz, 1H), 2.06 (d, J=14.2Hz, 1H), 2.43 (d, J=13.4Hz, 1H), 2.60 (d, J=13.4Hz, 1H), 5.69 (s, 1H), 7.00 (s, 2H), 7.31 (td, J=7.3Hz, 1.2Hz, 1H), 7.34-7.42 (m, 2H), 7.45 (d, J=8.1Hz, 1H), 7.55-7.61 (m, 2H), 7.66 (d, J=8.1Hz, 1H), 7.88 (d, J=16.4Hz, 1H), 7.96-8.03 (m, 3H), 8.26 (d, J=7.3Hz, 1H), 8.39 (d, J=8.8Hz, 1H).

Reaction was carried out in the same manner as in
 30 Examples 1 to 3 to obtain compounds of the following
 Example 17.

Example 17

Sodium [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-yl]oxy-
 35 acetate (sodium salt of Exemplary compound 505)

Appearance; pale yellowish solid

¹H-NMR (δ, DMSO-d₆); 3.79 (bs, 2H), 5.08 (bs, 1H), 7.06-7.35 (m, 3H), 7.40-7.55 (m, 4H), 7.62-7.72 (m, 1H), 7.92-7.97 (m, 1H), 7.93 (d, J=16.4Hz, 1H), 8.00 (d, J=9.8Hz, 1H), 8.01 (d, J=8.5Hz, 1H), 8.23 (d, J=7.3Hz, 1H), 8.38 (d, J=8.8Hz, 1H).

Reaction was carried out in the same manner as in Example 4 to obtain a compound of the following Example 18.

Sodium 3-([3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-yl]thio)-propionate (sodium salt of Exemplary compound 510)

Appearance; pale yellowish solid

¹H-NMR (δ, DMSO-d₆); 1.96-2.01 (m, 2H), 2.36-2.42 (m, 2H), 5.56 (s, 1H), 7.06-7.20 (m, 3H), 7.33 (d, J=6.6Hz, 1H), 7.41 (td, J=13.4Hz, 5.5Hz, 1H), 7.47 (d, J=8.1Hz, 1H), 7.55 (d, J=16.4Hz, 1H), 7.67 (dd, J=9.4Hz, 1.5Hz, 1H), 7.83-7.89 (m, 2H), 7.96 (d, J=8.6Hz, 1H), 8.01 (d, J=9.7Hz, 1H), 8.21 (d, J=7.3Hz, 1H), 8.38 (d, J=8.8Hz, 1H).

Reaction was carried out in the same manner as in Examples 1 to 3 to obtain a compound of the following Example 19.

Example 19

Sodium [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-9-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetate (sodium salt of Exemplary compound 535)

Appearance; pale yellowish solid

¹H-NMR (δ, DMSO-d₆); 3.79 (s, 2H), 5.07 (s, 1H), 7.35-7.51 (m, 6H), 7.66 (bs, 2H), 7.90-8.03 (m, 4H), 8.23 (d, J=7.3 Hz, 1H), 8.38 (d, J=8.5Hz, 1H).

Reaction was carried out in the same manner as in Example 4 to obtain compounds of the following Examples 20 to 21.

Example 20

Sodium 3-([3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-9-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thio)-propionate (sodium salt of Exemplary compound 542)

Appearance; yellowish solid

¹H-NMR (δ, DMSO-d₆); 1.97-2.09 (m, 2H), 2.41-2.46 (m, 2H),
5.56 (s, 1H), 7.20 (d, J=12.5Hz, 1H), 7.26 (d, J=12.5Hz,
1H), 7.36-7.52 (m, 4H), 7.56 (d, J=16.4Hz, 1H), 7.69 (dd,
5 J=8.1Hz, 1.5Hz, 1H), 7.88 (d, J=16.4Hz, 1H), 7.91 (d,
J=1.5Hz, 1H), 7.98 (d, J=8.8Hz, 1H), 8.02 (d, J=9.8Hz, 1H),
8.23 (d, J=7.3Hz, 1H), 8.40 (d, J=8.6Hz, 1H).

Example 21

Sodium 3-{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-
10 yl)ethenyl]-8-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thio}-
propionate (sodium salt of Exemplary compound 549)

Appearance; yellowish solid

¹H-NMR (δ, DMSO-d₆); 1.96-2.01 (m, 2H), 2.35-2.40 (m, 2H),
5.54 (s, 1H), 6.98 (d, J=12.0Hz, 1H), 7.07 (d, J=12.2Hz,
15 1H), 7.43 (dd, J=8.3Hz, 2.2Hz, 1H), 7.47 (d, J=8.1Hz, 1H),
7.50-7.58 (m, 2H), 6.68 (d, J=7.8Hz, 1H), 7.84-7.90 (m,
2H), 7.98 (d, J=8.8Hz, 1H), 8.02 (d, J=9.8Hz, 1H), 8.02 (d,
J=7.8Hz, 1H), 8.39 (d, J=8.8Hz, 1H).

Reaction was carried out in the same manner as in
20 Examples 1 to 3 to obtain compounds of the following
Examples 22 to 25.

Example 22

Sodium [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-
ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]oxy-
25 acetate (sodium salt of Exemplary compound 557)

Appearance; pale yellowish solid

¹H-NMR (δ, DMSO-d₆); 3.06 (d, J=15.0Hz, 1H), 3.12 (d,
J=15.0Hz, 1H), 6.52 (s, 1H), 7.07 (d, J=12.0Hz, 1H), 7.11
(d, J=12.5Hz, 1H), 7.36 (t, J=7.8Hz, 1H), 7.48-7.58 (m,
30 4H), 7.78-7.82 (m, 2H), 7.94 (d, J=16.4Hz, 1H), 7.99-8.02
(m, 2H), 8.22 (d, J=7.3Hz, 1H), 8.39 (d, J=8.8Hz, 1H).

Example 23

Sodium 2-(R)-{[3-[(E)-2-(7-chloro-6-fluoroquinolin-
2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}propion-
35 ate (sodium salt of Exemplary compound 558)

Appearance; yellowish solid

¹H-NMR (δ, DMSO-d₆); 0.84-0.86, 1.46-1.50 (each m, 1H in total), 3.00-3.04, 3.75-3.79 (each m, 1H in total), 5.01, 5.95 (each s, 1H in total), 7.03, 7.15 (each s, 1H in total), 7.20-8.01 (m, 11H), 8.04 (d, J=10.0Hz, 1H), 8.23 (d, J=7.3Hz, 1H), 8.37 (dd, J=8.8Hz, 2.4Hz, 1H).

Example 24

Sodium 2-(S)-{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}propionate (sodium salt of Exemplary compound 558)

10 Appearance; pale yellowish solid

¹H-NMR (δ, DMSO-d₆); 0.82-0.84, 1.44-1.48 (each m, 1H in total), 3.33-3.42, 3.69-3.78 (each m, 1H in total), 5.00, 5.94 (each s, 1H in total), 7.03, 7.15 (each s, 1H in total), 7.20-7.99 (m, 11H), 8.04 (d, J=10.0Hz, 1H), 8.23 (d, J=7.3Hz, 1H), 8.37 (dd, J=8.8Hz, 2.4Hz, 1H).

Example 25

Sodium 3-{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}-propionate (sodium salt of Exemplary compound 559)

20 Appearance; pale yellowish solid

¹H-NMR (δ, DMSO-d₆); 1.86-1.92 (m, 2H), 3.09-3.22 (m, 2H), 6.26 (s, 1H), 7.08 (d, J=12.0Hz, 1H), 7.13 (d, J=12.2Hz, 1H), 7.38 (t, J=7.6Hz, 1H), 7.51 (dd, J=7.8Hz, 1.2Hz, 1H), 7.53-7.62 (m, 3H), 7.80 (dd, J=8.1Hz, 1.5Hz, 1H), 7.92 (s, 1H), 7.95 (d, J=16.4Hz, 1H), 7.99-8.03 (m, 2H), 8.22 (d, J=7.3Hz, 1H), 8.39 (d, J=8.5Hz, 1H).

Reaction was carried out in the same manner as in Example 4 to obtain compounds of the following Examples 26 to 28.

30 Example 26

Sodium [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thioacetate (sodium salt of Exemplary compound 562)

Appearance; orange solid

35 ¹H-NMR (δ, DMSO-d₆); 2.77 (d, J=13.4Hz, 1H), 2.85 (d, J=13.7Hz, 1H), 6.17 (s, 1H), 7.04 (d, J=12.0Hz, 1H), 7.08

(d, J=12.0Hz, 1H), 7.30 (t, J=7.8Hz, 1H), 7.41 (d, J=7.1Hz, 1H), 7.47-7.58 (m, 3H), 7.72 (d, J=8.1Hz, 1H), 7.79 (s, 1H), 7.91 (d, J=16.4Hz, 1H), 7.98-8.03 (m, 2H), 8.22 (d, J=7.7Hz, 1H), 8.39 (d, J=8.8Hz, 1H).

5 Example 27

Sodium 3-{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thio}-propionate (sodium salt of Exemplary compound 564)

Appearance; yellowish solid

10 ¹H-NMR (δ, DMSO-d₆); 1.93-2.07 (m, 2H), 2.49-2.55 (m, 2H), 6.02 (s, 1H), 7.06 (d, J=12.2Hz, 1H), 7.11 (d, J=12.0Hz, 1H), 7.32 (t, J=7.6Hz, 1H), 7.42 (dd, J=7.8Hz, 1.2Hz, 1H), 7.51 (d, J=8.1Hz, 1H), 7.52-7.62 (m, 2H), 7.75 (dd, J=8.1Hz, 1.5Hz, 1H), 7.82 (d, J=1.5Hz, 1H), 7.92 (d, J=16.4Hz, 1H), 7.99-8.03 (m, 2H), 8.22 (d, J=7.3Hz, 1H), 15 7.39 (d, J=8.8Hz, 1H).

Example 28

Sodium 3-{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionate (sodium salt of Exemplary compound 613)

20 Appearance; pale yellowish solid

FAB-MS (m/z): 492 (M⁺+1).

¹H-NMR (δ, DMSO-d₆); 1.99-2.09 (m, 2H), 2.43-2.53 (m, 2H), 5.65 (s, 1H), 7.69-7.30 (m, 2H), 7.50-7.59 (m, 3H), 7.67- 25 7.74 (m, 2H), 7.85-8.03 (m, 5H), 8.23 (d, J=7.3Hz, 1H), 8.40 (d, J=8.8Hz, 1H).

Reaction was carried out in the same manner as in Examples 1 to 3 to obtain a compound of the following Example 29.

30 Example 29

Sodium [3-[(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetate (sodium salt of Exemplary compound 1411)

Appearance; orange solid

35 ¹H-NMR (δ, DMSO-d₆); 3.80 (bs, 2H), 5.04 (bs, 1H), 7.16 (bs, 2H), 7.27-7.66 (m, 8H), 7.92-8.02 (m, 4H), 8.04 (d,

J=2.2Hz, 1H), 8.40 (d, J=8.5Hz, 1H).

Reaction was carried out in the same manner as in Example 4 to obtain compounds of the following Examples 30 to 31.

5 Example 30

Sodium [3-[(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thioacetate (sodium salt of Exemplary compound 1416)

Appearance; pale yellowish solid

10 FAB-MS (m/z); 492 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 2.65 (d, J=13.9Hz, 2H), 2.68 (d, J=13.9Hz, 2H), 5.64 (s, 1H), 6.98 (s, 2H), 7.22-7.46 (m, 5H), 7.53 (d, J=16.4Hz, 1H), 7.59 (dd, J=8.8Hz, 2.2Hz, 1H), 7.67 (dd, J=8.1Hz, 1.5Hz, 1H), 7.77 (d, J=1.5Hz, 1H), 7.87-
15 8.04 (m, 4H), 8.41 (d, J=8.5Hz, 1H).

Example 31

Sodium 3-{[3-[(E)-2-(5,6,7,8-tetrahydroquinolin-2-yl)ethenyl]-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionate (sodium salt of Exemplary compound 1921)

20 Appearance; pale yellowish solid

FAB-MS (m/z): 544 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 1.75-1.85 (m, 4H), 2.00-2.05 (m, 2H), 2.42-2.47 (m, 2H), 2.72-2.87 (m, 4H), 5.60 (s, 1H), 7.14-7.32 (m, 4H), 7.44-7.47 (m, 2H), 7.53-7.61 (m, 3H), 7.72
25 (d, J=7.6Hz, 1H), 7.81-7.83 (m, 2H).

Reference example 1

(a) 7-Chloro-2-vinyl quinolin

The title compound was obtained by the method as described in J. Org. Chem., 61, 3398 (1996).

30 Appearance; ocherous solid

$^1\text{H-NMR}$ (δ , DMSO- d_6); 5.69 (dd, J=11.0Hz, 1.0Hz, 1H), 6.31 (dd, J=17.6Hz, 0.7Hz, 1H), 7.00 (dd, J=17.8Hz, 11.0Hz, 1H), 7.45 (dd, J=8.8Hz, 2.0Hz, 1H), 7.60 (d, J=8.5Hz, 1H), 7.71 (d, J=8.5Hz, 1H), 8.06-8.10 (m, 2H).

35 (b) 3-Trifluoromethanesulfonyloxy-5H-dibenzo[a,d]-cyclohepten-5-one

In 30 ml of methylene chloride was dissolved 0.23 g (1.0 mmol) of 3-hydroxy-5H-dibenzo[a,d]cyclohepten-5-one and after cooling the solution with ice-water, 0.33 ml (2.0 mmol) of trifluoromethanesulfonic anhydride and 0.27 ml (2.0 mmol) of triethylamine were added to the solution and the resulting mixture was stirred under ice-cooling for 4 hours.

After completion of the reaction, the reaction mixture was concentrated, and the residue was applied to silica gel column chromatography (eluent: chloroform) to obtain 0.30 g of the title compound as brown oily product. CI-MS (m/z); 355 ($M^+ + 1$).

EI-MS; 354 (M^+).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 7.31 (d, $J=12.2\text{Hz}$, 1H), 7.37 (d, $J=12.2\text{Hz}$, 1H), 7.50-8.40 (m, 7H).

(c) 3-[(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-one

In 30 ml of N,N-dimethylformamide were dissolved 0.89 g (2.5 mmol) of 3-trifluoromethanesulfonyloxy-5H-dibenzo[a,d]cyclohepten-5-one and 0.48 g (2.5 mmol) of 7-chloro-2-vinyl quinoline, and then, 100 mg (0.5 mmol) of palladium acetate, 420 mg (1.6 mmol) of triphenylphosphine and 2.1 g (24 mmol) of lithium bromide were added to the mixture and the atmosphere thereof was made nitrogen. Then, 5.0 ml (36 mmol) of triethylamine was added to the mixture, and the resulting mixture was stirred under nitrogen atmosphere at 120°C for 10 hours.

After completion of the reaction, the reaction mixture was concentrated, and the residue was applied to silica gel column chromatography (eluent: toluene/ethyl acetate=9/1 (volume ratio)) to obtain 0.53 g of the title compound as pale brown solid.

CI-MS (m/z); 394 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 7.28 (s, 2H), 7.59-7.69 (m, 3H), 7.78-7.83 (m, 3H), 7.97-8.06 (m, 4H), 8.13-8.20 (m, 2H), 8.38 (s, 1H), 8.44 (d, $J=9.3\text{Hz}$, 1H).

(d) 3-[(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-ol

In a mixed solution of 40 ml of tetrahydrofuran and 10 ml of methanol was dissolved 0.53 g (1.3 mmol) of 3-
5 [(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]-cyclohepten-5-one, and 0.11 g (2.6 mmol) of sodium borohydride was added to the mixture and the resulting mixture was stirred at room temperature for 3 hours.

After completion of the reaction, water was added to
10 the reaction mixture, a pH of the mixture was adjusted to about 2.0 with 1N hydrochloric acid and the mixture was extracted with ethyl acetate. The organic layer was washed with a saturated aqueous sodium chloride solution, dried over anhydrous magnesium sulfate and concentrated. The
15 residue was dried under reduced pressure to obtain 0.33 g of the title compound as ochreous solid.

CI-MS (m/z); 396 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 5.10 (d, $J=3.4\text{Hz}$, 1H), 6.20 (s, 1H),
7.18 (s, 2H), 7.26 (td, $J=8.8\text{Hz}$, 1.2Hz, 1H), 7.37-7.52 (m,
20 4H), 7.59 (dd, $J=8.5\text{Hz}$, 2.2Hz, 1H), 7.64 (dd, $J=8.1\text{Hz}$, 1.7Hz, 1H), 7.76 (d, $J=7.6\text{Hz}$, 1H), 7.91-8.04 (m, 5H), 8.40 (d, $J=8.3\text{Hz}$, 1H).

Reference example 2

Reaction was carried out in the same manner as in
25 Reference example 1(a) to obtain a compound of the following Reference example 2(a).

(a) 7-Chloro-6-fluoro-2-vinylquinoline

Appearance; yellowish solid

CI-MS (m/z); 208 ($M^+ + 1$).

30 EI-MS; 207 (M^+).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 5.73 (dd, $J=11.0\text{Hz}$, 1.0Hz, 1H), 6.43 (dd, $J=17.6\text{Hz}$, 1.0Hz, 1H), 6.98 (dd, $J=17.8\text{Hz}$, 11.0Hz, 1H),
7.89 (d, $J=8.8\text{Hz}$, 1H), 8.01 (d, $J=9.8\text{Hz}$, 1H), 8.21 (d, $J=7.3\text{Hz}$, 1H). 8.37 (d, $J=9.0\text{Hz}$, 1H).

35 Reaction was carried out in the same manner as in Reference example 1(c) to obtain a compound of the follow-

ing Reference example 2(b).

(b) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-one

Appearance; pale yellowish solid

5 $^1\text{H-NMR}$ (δ , DMSO- d_6); 7.29 (s, 2H), 7.63-7.70 (m, 2H), 7.78-7.84 (m, 3H), 8.00-8.05 (m, 3H), 8.13-8.21 (m, 2H), 8.23 (d, $J=7.6\text{Hz}$, 1H), 8.38 (d, $J=1.7\text{Hz}$, 1H), 8.42 (d, $J=9.3\text{Hz}$, 1H).

10 Reaction was carried out in the same manner as in Reference example 1(d) to obtain a compound of the following Reference example 2(c).

(c) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-5H-dibenzo[a,d]cyclohepten-5-ol

Appearance; yellowish solid

15 $^1\text{H-NMR}$ (δ , DMSO- d_6); 5.09 (d, $J=3.9\text{Hz}$, 1H), 6.20 (s, 1H), 7.18 (s, 2H), 7.26 (td, $J=7.3\text{Hz}$, 1.2Hz, 1H), 7.37-7.50 (m, 4H), 7.60 (d, $J=7.6\text{Hz}$, 1H), 7.63 (dd, $J=7.8\text{Hz}$, 1.7Hz, 1H), 7.93 (d, $J=16.4\text{Hz}$, 1H), 7.98-8.02 (m, 3H), 8.21 (d, $J=7.3\text{Hz}$, 1H), 8.38 (d, $J=8.5\text{Hz}$, 1H).

20 Reference example 3

Reaction was carried out in the same manner as in Reference example 1(a) to obtain a compound of the following Reference example 3(a).

(a) 6,7-Difluoro-2-vinylquinoline

25 Appearance; pale yellowish solid

CI-MS (m/z); 192 (M^++1).

EI-MS; 191 (M^+).

30 $^1\text{H-NMR}$ (δ , DMSO- d_6); 5.69 (dd, $J=11.0\text{Hz}$, 1.0Hz, 1H), 6.39 (dd, $J=17.6\text{Hz}$, 1.0Hz, 1H), 6.93 (dd, $J=8.1\text{Hz}$, 3.9Hz, 1H), 6.95 (dd, $J=17.6\text{Hz}$, 11.0Hz, 1H), 7.82 (d, $J=8.8\text{Hz}$, 1H), 8.00 (dd, $J=9.0\text{Hz}$, 2.2Hz, 1H), 8.33 (d, $J=8.5\text{Hz}$, 1H).

Reaction was carried out in the same manner as in Reference example 1(c) to obtain a compound of the following Reference example 3(b).

35 (b) 3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-one

Appearance; brown solid

CI-MS (m/z); 396 ($M^+ + 1$).

EI-MS; 395 (M^+).

^1H -NMR (δ , DMSO- d_6); 7.29 (s, 2H), 7.50-7.90 (m, 7H), 7.95-8.30 (m, 5H), 8.41 (d, $J=8.5\text{Hz}$, 1H).

Reaction was carried out in the same manner as in Reference example 1(d) to obtain a compound of the following Reference example 3(c).

(c) 3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-ol

Appearance; brown solid

CI-MS (m/z); 398 ($M^+ + 1$).

EI-MS; 397 (M^+).

^1H -NMR (δ , DMSO- d_6); 5.09 (s, 1H), 7.18 (s, 2H), 7.20-7.70 (m, 2H), 7.28 (d, $J=7.3\text{Hz}$, 1H), 7.38 (d, $J=7.6\text{Hz}$, 1H), 7.43 (d, $J=7.8\text{Hz}$, 1H), 7.49 (d, $J=7.6\text{Hz}$, 1H), 7.76 (d, $J=7.6\text{Hz}$, 1H), 7.80-8.20 (m, 4H), 7.93 (d, $J=16.6\text{Hz}$, 1H), 8.36 (d, $J=8.5\text{Hz}$, 1H).

Reference example 4

(a) 5,6,7,8-Tetrahydro-2-vinylquinoline

In 300 ml of acetonitrile was dissolved 14.7 g (65 mmol) of (5,6,7,8-tetrahydroquinolin-2-yl)methyl bromide, and 25.5 g (97.5 mmol) of triphenylphosphine was added to the mixture and the resulting mixture was refluxed for 2 hours. After cooling the mixture with an ice-bath, precipitates were collected by filtration and washed with diethyl ether to obtain (5,6,7,8-tetrahydroquinolin-2-yl)methyltriphenylphosphonium bromide as pale yellowish solid. Then, the obtained solid was dissolved in 250 ml of chloroform, 9.50 g (117 mmol) of 37% formaldehyde and 8.10 g (76 mmol) of an aqueous sodium carbonate solution (20 ml) were added to the solution and the mixture was stirred at room temperature for 2 hours.

After completion of the reaction, water was added to the reaction mixture and the mixture was extracted with chloroform. The organic layer was dried over potassium

carbonate and concentrated. The residue was applied to silica gel column chromatography (eluent: hexane/ethyl acetate=9/1 (volume ratio)) to obtain 6.01 g of the title compound as yellowish oily product.

5 CI-MS (m/z); 160 ($M^+ + 1$).

EI-MS; 159 (M^+).

$^1\text{H-NMR}$ (δ , CDCl_3); 1.85 (m, 4H), 2.75 (dd, $J=17.6\text{Hz}$, 1.0Hz , 2H), 2.91 (dd, $J=8.1\text{Hz}$, 3.9Hz , 2H), 5.41 (dd, $J=10.7\text{Hz}$, 1.2Hz , 1H), 6.06 (dd, $J=17.6\text{Hz}$, 1.2Hz , 1H), 6.75 (dd, $J=17.6\text{Hz}$, 10.7Hz , 1H), 7.14 (d, $J=7.8\text{Hz}$, 1H), 7.31 (d, $J=7.8\text{Hz}$, 1H).

Reaction was carried out in the same manner as in Reference example 1(c) to obtain a compound of the following Reference example 4(b).

15 (b) 3-[(E)-2-(5,6,7,8-tetrahydroquinolin-2-yl)-ethenyl]-5H-dibenzo[a,d]cyclohepten-5-one

Appearance; pale yellowish solid

$^1\text{H-NMR}$ (δ , CDCl_3); 1.81-1.95 (m, 4H), 2.79 (t, $J=6.1$, 2H), 2.96 (t, $J=6.4\text{Hz}$, 2H), 7.05 (s, 2H), 7.15-7.37 (m, 4H), 7.52-7.64 (m, 4H), 7.81-7.85 (m, 1H), 8.23-8.26 (m, 1H), 8.40 (d, $J=2.0\text{Hz}$, 1H).

Reaction was carried out in the same manner as in Reference example 1(d) to obtain a compound of the following Reference example 4(c).

25 (c) 3-[(E)-2-(5,6,7,8-tetrahydroquinolin-2-yl)-ethenyl]-5H-dibenzo[a,d]cyclohepten-5-ol

Appearance; pale yellowish solid

$^1\text{H-NMR}$ (δ , CDCl_3); 1.80-2.00 (m, 4H), 2.77 (m, 2H), 2.95 (m, 2H), 5.45 (s, 1H), 7.09 (s, 1H), 7.20-7.70 (m, 10H), 7.89 (s, 1H).

Reference example 5

(a) 3-(7-chloro-6-fluoroquinolin-2-yl)methoxy-5H-dibenzo[a,d]cyclohepten-5-one

In 30 ml of N,N-dimethylformamide were dissolved 0.37 g (1.4 mmol) of 2-bromomethyl-7-chloro-6-fluoro-quinoline and 0.30 g (1.4 mmol) of 3-hydroxy-5H-dibenzo-

[a,d]cyclohepten-5-one, and 0.37 g (2.8 mmol) of potassium carbonate was added to the mixture and the resulting mixture was stirred at room temperature for 6 hours.

After completion of the reaction, water was added to the reaction mixture and the mixture was extracted with ethyl acetate. The organic layer was washed with a saturated aqueous sodium chloride solution, dried over anhydrous magnesium sulfate and concentrated. The residue was dried under reduced pressure to obtain 0.47 g of the title compound as dark brown solid.

CI-MS (m/z); 416 ($M^+ + 1$).

EI-MS; 415 (M^+).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 5.53 (s, 2H), 7.13 (d, $J=12.2\text{Hz}$, 1H), 7.21 (d, $J=12.2\text{Hz}$, 1H), 7.51 (dd, $J=8.5\text{Hz}$, 2.7Hz, 1H), 7.55-7.90 (m, 6H), 8.00-8.35 (m, 2H), 8.31 (d, $J=7.3\text{Hz}$, 1H), 8.46 (d, $J=8.5\text{Hz}$, 1H).

(b) 3-(7-chloro-6-fluoroquinolin-2-yl)methoxy-5H-dibenzo[a,d]cyclohepten-5-ol

Reaction was carried out in the same manner as in Reference example 1(d) except for using 3-(7-chloro-6-fluoroquinolin-2-yl)methoxy-5H-dibenzo[a,d]cyclohepten-5-one in place of 3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-one to obtain the title compound.

Appearance; brown solid

CI-MS (m/z); 418 ($M^+ + 1$).

EI-MS; 418 (M^+).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 5.38 (s, 1H), 5.41 (s, 2H), 6.80-7.90 (m, 9H), 7.95-8.35 (m, 3H), 8.42 (d, $J=8.8\text{Hz}$, 1H).

Reference example 6

(a) 2-methoxycarbonylbenzyltriphenylphosphonium bromide

In 300 ml of acetonitrile were dissolved 86.32 g (377 mmol) of methyl 2-(bromomethyl)benzoate and 98.84 g (377 mmol) of triphenylphosphine, and the solution was refluxed for 2 hours.

After completion of the reaction, 700 ml of diethyl ether was added to the reaction mixture. The precipitated white precipitates were collected by filtration and washed with diethyl ether. The precipitates were dried under
5 reduced pressure to obtain 117.42 g of the title compound as pale brown solid.

CI-MS (m/z); 410 (M^+ -Br).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 3.46 (s, 3H), 5.53 (d, $J=15.4\text{Hz}$, 2H),
7.30-7.40 (m, 1H), 7.45-7.63 (m, 8H), 7.65-7.80 (m, 6H),
10 7.82-7.95 (m, 4H).

(b) Methyl 2-[2-(4-methoxyphenyl)ethenyl]benzoate

In 50 ml of acetonitrile were dissolved 4.91 g (10 mmol) of 2-methoxycarbonylbenzyltriphenylphosphonium bromide and 1.77 g (13 mmol) of 4-methoxybenzaldehyde, and
15 2.48 g (20 mmol) of 1,5-diazabicyclo[4.3.0]-5-nonene was added to the solution and the mixture was refluxed for 5 hours.

After completion of the reaction, the reaction mixture was concentrated, and the residue was applied to
20 silica gel column chromatography (eluent: toluene) to obtain 1.25 g of the title compound (E,Z mixture) as pale yellowish oily product.

CI-MS (m/z); 269 (M^++1).

EI-MS; 268 (M^+).

25 Cis isomer (Z isomer)

$^1\text{H-NMR}$ (δ , DMSO- d_6); 3.79 (s, 3H), 3.87 (s, 3H), 6.74 (d, $J=9.0\text{Hz}$, 1H), 6.90-7.00 (m, 2H), 7.10-7.73 (m, 5H), 7.75-7.95 (m, 2H).

Trans isomer (E isomer)

30 $^1\text{H-NMR}$ (δ , DMSO- d_6); 3.69 (s, 3H), 3.87 (s, 3H), 6.56 (d, $J=12.2\text{Hz}$, 1H), 6.74 (d, $J=9.0\text{Hz}$, 1H), 6.84 (d, $J=12.2\text{Hz}$, 1H), 6.90-7.00 (m, 2H), 7.10-7.73 (m, 3H), 7.75-7.95 (m, 2H).

(c) 2-[2-(4-Methoxyphenyl)ethenyl]benzoic acid

35 To 2 ml of a methanol solution containing 1.20 g (4.5 mmol) of methyl 2-[2-(4-methoxyphenyl)ethenyl]benzoic

acid was added 1.5 g (26.7 mmol) of an aqueous potassium hydroxide solution (20 ml), and the mixture was refluxed for 6 hours.

After completion of the reaction, the reaction mixture was cooled and a 1N aqueous hydrochloric acid solution was added to the mixture to adjust a pH thereof to 3. Formed white precipitates were collected by filtration, washed with water and dried under reduced pressure to obtain 1.12 g of the title compound (E,Z mixture) as white solid.

CI-MS (m/z); 255 ($M^+ + 1$).

EI-MS; 254 (M^+).

Cis isomer (Z isomer)

$^1\text{H-NMR}$ (δ , DMSO- d_6); 3.78 (s, 3H), 6.53 (d, $J=12.2\text{Hz}$, 1H), 6.74 (d, $J=8.8\text{Hz}$, 1H), 6.89 (d, $J=12.2\text{Hz}$, 1H), 6.96 (d, $J=8.8\text{Hz}$, 1H), 6.97 (d, $J=8.8\text{Hz}$, 1H), 7.15-7.60 (m, 3H), 7.70-7.95 (m, 2H), 12.93 (br, 1H).

Trans isomer (E isomer)

$^1\text{H-NMR}$ (δ , DMSO- d_6); 3.69 (s, 3H), 6.74 (d, $J=8.8\text{Hz}$, 1H), 6.96 (d, $J=8.8\text{Hz}$, 1H), 6.97 (d, $J=8.8\text{Hz}$, 1H), 7.12 (d, $J=16.1\text{Hz}$, 1H), 7.15-7.60 (m, 4H), 7.70-7.95 (m, 2H), 12.93 (br, 1H).

(d) 2-[2-(4-Methoxyphenyl)ethyl]benzoic acid

To 20 ml of a methanol solution containing 0.20 g (1.3 mmol) of 2-[2-(4-methoxyphenyl)ethenyl]benzoic acid was added 0.10 g of 10% palladium-activated charcoal, and then, the mixture was stirred at room temperature for 5 hours under hydrogen atmosphere (normal pressure).

After completion of the reaction, 10% palladium-activated charcoal was removed from the reaction mixture by filtration by using Celite (tradename), and the filtrate was concentrated to obtain 0.18 g of the title compound as pale yellowish solid.

CI-MS (m/z); 257 ($M^+ + 1$).

EI-MS; 256 (M^+).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 2.70-2.80 (m, 2H), 3.10-3.20 (m, 2H),

3.67 (s, 3H), 6.80-6.90 (m, 2H), 7.10-7.20 (m, 2H), 7.22-7.35 (m, 2H), 7.40-7.50 (m, 1H), 7.75-7.85 (m, 1H), 12.88 (br, 1H).

5 (e) 3-Methoxy-10,11-dihydro-5H-dibenzo[a,d]cyclohepten-5-one

To 50 ml of a methylene chloride solution containing 1.80 g (7.0 mmol) of 2-[2-(4-methoxyphenyl)ethyl]benzoic acid were added 1.83 ml (13 mmol) of trifluoroacetic anhydride and 0.8 ml (6.3 mmol) of boron trifluoride ethyl ether complex and the mixture was stirred at room temperature for 5 hours.

After completion of the reaction, the reaction mixture was concentrated, and the residue was applied to silica gel column chromatography (eluent: toluene) to obtain 0.65 g of the title compound as yellow oily product. CI-MS (m/z); 239 ($M^+ + 1$).

EI-MS; 238 (M^+).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 3.12 (s, 4H), 3.79 (s, 3H), 7.12 (dd, $J=8.3\text{Hz}$, 2.9Hz , 1H), 7.28 (d, $J=8.3\text{Hz}$, 1H), 7.35 (d, $J=7.8\text{Hz}$, 2H), 7.40 (d, $J=2.9\text{Hz}$, 1H), 7.51 (dd, $J=7.3\text{Hz}$, 1.5Hz , 1H), 7.84 (d, $J=8.1\text{Hz}$, 1H).

(f) 3-Methoxy-5H-dibenzo[a,d]cyclohepten-5-one

To 20 ml of a 1,2-dichloroethane solution containing 0.65 g (2.7 mmol) of 3-methoxy-10,11-dihydro-5H-dibenzo[a,d]cyclohepten-5-one were added 0.80 g (4.5 mmol) of N-bromosuccineimide and 0.14 g (0.6 mmol) of benzoyl peroxide, and the mixture was refluxed for 3 hours.

After completion of the reaction, 100 ml of chloroform was added to the reaction mixture and the resulting mixture was washed with 50 ml of an aqueous 1N sodium hydroxide solution. The organic layer was washed with a saturated aqueous sodium chloride solution, dried over anhydrous magnesium sulfate and concentrated. The residue was dissolved in 50 ml of 1,2-dichloroethane, 50 ml of triethylamine was added thereto and the resulting mixture was refluxed for 2 hours.

After completion of the reaction, the reaction mixture was concentrated, and the residue was applied to silica gel column chromatography (eluent: toluene/ethyl acetate=9/1 (volume ratio)) to obtain 0.57 g of the title
5 compound as brown solid.

CI-MS (m/z); 237 ($M^+ + 1$).

EI-MS; 236 (M^+).

^1H -NMR (δ , DMSO- d_6); 3.90 (s, 3H), 7.12 (d, $J=12.2\text{Hz}$, 1H),
7.22 (d, $J=12.2\text{Hz}$, 1H), 7.38 (dd, $J=8.5\text{Hz}$, 2.9Hz , 1H),
10 7.55-7.68 (m, 2H), 7.70-7.80 (m, 3H), 8.13 (d, $J=7.6\text{Hz}$, 1H).

(g) 3-Hydroxy-5H-dibenzo[a,d]cyclohepten-5-one

In 100 ml of acetic acid was dissolved 2.03 g (8.6 mmol) of 3-methoxy-5H-dibenzo[a,d]cyclohepten-5-one, 20 ml
15 of 48% hydrobromic acid was added to the solution and the resulting mixture was refluxed for 10 hours.

After completion of the reaction, the reaction mixture was poured into 200 ml of ice-water and the formed solid was collected by filtration, washed with water and
20 dried to obtain 1.48 g of the title compound as dark brown solid.

CI-MS (m/z); 223 ($M^+ + 1$).

EI-MS; 222 (M^+).

^1H -NMR (δ , DMSO- d_6); 7.04 (d, $J=12.0\text{Hz}$, 1H), 7.12 (d,
25 $J=12.0\text{Hz}$, 1H), 7.19 (dd, $J=8.5\text{Hz}$, 2.9Hz , 1H), 7.50-7.64 (m, 3H), 7.67-7.80 (m, 2H), 8.11 (d, $J=7.6\text{Hz}$, 1H).

Reference example 7

(a) 3-(4-Methoxybenzylidene)phthalide

22.28 g (150 mmol) of phthalic anhydride, 25.0 g
30 (150 mmol) of 4-methoxyphenyl acetate and 2.0 g (24 mmol) of sodium acetate were mixed and the mixture was stirred at 220°C for 8 hours.

After completion of the reaction, ethanol was added to the reaction mixture and insoluble material was removed
35 by filtration, and the filtrate was concentrated. The obtained residue was applied to silica gel column chromato-

graphy (eluent: toluene/ethyl acetate=9/1 (volume ratio)) to obtain 13.15 g of the title compound as yellowish solid. CI-MS (m/z); 253 ($M^+ + 1$).

EI-MS; 252 (M^+).

- 5 $^1\text{H-NMR}$ (δ , DMSO-d_6); 3.81 (s, 3H), 6.90 (s, 1H), 7.06 (d, $J=8.8\text{Hz}$, 1H), 7.64 (t, $J=7.3\text{Hz}$, 1H), 7.78 (d, $J=8.8\text{Hz}$, 1H), 7.87 (t, $J=7.3\text{Hz}$, 1H), 7.95 (d, $J=7.8\text{Hz}$, 1H), 8.09 (d, $J=7.8\text{Hz}$, 1H).

(b) 2-[2-(4-Methoxyphenyl)ethyl]benzoic acid

- 10 To 280 ml of an ethanol solution containing 20.0 g (79.3 mmol) of 3-(4-methoxybenzylidene)phthalide were added 16.6 ml (119 mmol) of triethylamine and Raney nickel (available from Kawaken Finechemical, developed nickel catalyst NTD-65, 24 ml), and the resulting mixture was stirred at 80°C
15 for 2.5 hours under pressure with hydrogen(10 atm).

- After completion of the reaction, the catalyst was removed from the reaction mixture by filtration using Celite (tradename), and the filtrate was concentrated. The obtained residue was dissolved in 200 ml of methylene
20 chloride, washed successively with 100 ml of a 1N aqueous hydrochloric acid solution, and a saturated aqueous sodium chloride solution, dried over anhydrous magnesium sulfate and concentrated. The residue was dried under reduced pressure to obtain 17.2 g of the title compound as white
25 solid. The obtained compound was the same as the compound of Reference example 6(d).

Reference example 8

(a) 2-[2-(4-Bromophenyl)ethenyl]-3-fluorobenzoic acid

- 30 To a liquor in which 11.6 g (121 mmol) of sodium t-butoxide had been suspended in 200 ml of N,N-dimethylformamide were added dropwise a solution in which 20.2 g (120 mmol) of methyl 3-fluoro-2-methylbenzoate and 22.3 g (120 mmol) of p-bromobenzaldehyde had been dissolved in 100 ml
35 of dimethylformamide at 5°C over 1 hour. After completion of dropwise addition, the mixture was stirred at 10°C for 1

hour, and further stirred at at room temperature overnight.

After completion of the reaction, 1.0 liter of water was added to the reaction mixture and the resulting mixture
5 was adjusted with conc. hydrochloric acid to pH 3.

Precipitated solid was collected by filtration, washed with water and dried under reduced pressure to obtain 25.3 g of the title compound as pale yellowish solid.

¹H-NMR (δ, CDCl₃); 7.06 (dd, J=16.9Hz, 1.7Hz, 1H), 7.38-
10 7.61 (m, 7H), 7.67 (dd, J=7.1Hz, 1.7Hz, 1H), 13.28 (b, 1H).

(b) 2-[2-(4-Bromophenyl)ethyl]-3-fluorobenzoic acid

To 25.3 g (78.8 mmol) of 2-[2-(4-bromophenyl)ethen-
yl]-3-fluorobenzoic acid were added 6.2 g (199.4 mmol) of
red phosphorus, 60 ml of 57% hydroiodic acid and 150 ml of
15 acetic acid, and the mixture was refluxed for 10 hours.
After cooling by allowing to stand, red phosphorus was
further added to the mixture and the mixture was refluxed.
Addition of red phosphorus was carried out further two
times.

20 After completion of the reaction, 1.0 liter of water
was added to the reaction mixture, precipitated solid was
collected by filtration, water was added to the solid and
the mixture was adjusted to pH 12 with an aqueous sodium
hydroxide solution, and insoluble material was filtered
25 off. The filtrate was adjusted to pH 3 with conc. hydro-
chloric acid, precipitated solid was collected by filtra-
tion, washed with water and dried under reduced pressure to
obtain 23.9 g of the title compound as colorless solid.

¹H-NMR (δ, CDCl₃); 2.76-2.81 (m, 2H), 3.15-3.20 (m, 2H),
30 7.16 (d, J=8.3Hz, 2H), 7.35-7.39 (m, 2H), 7.47 (d, J=8.3Hz,
2H), 7.65-7.68 (m, 1H), 13.4 (b, 1H).

(c) 3-Bromo-9-fluoro-10,11-dihydro-5H-dibenzo[a,d]-
cyclohepten-5-one

To 25.8 g (80 mmol) of 2-[2-(4-bromophenyl)ethyl]-3-
35 fluorobenzoic acid was added 550 g of polyphosphoric acid,
and the mixture was stirred at 170°C for 3 hours. After

cooling the mixture to 80°C by allowing to stand, water was added to the mixture and the resulting mixture was cooled to room temperature under stirring by allowing to stand.

After completion of the reaction, the reaction mixture was extracted with chloroform. The extract was washed successively with water, an aqueous sodium hydroxide solution, and water. The organic layer was dried over anhydrous sodium sulfate, concentrated and the residue was applied to silica gel column chromatography (eluent: hexane/ethyl acetate=9/1(volume ratio)) to obtain 18.6 g of the title compound as brown solid.

CI-MS (m/z); 305 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , CDCl_3); 3.11-3.23 (m, 4H), 7.17 (d, $J=8.5\text{Hz}$, 1H), 7.19-7.34 (m, 2H), 7.55 (dd, $J=8.3\text{Hz}$, 2.2Hz , 1H), 7.79 (d, $J=7.1\text{Hz}$, 1H), 7.99 (m, $J=2.2\text{Hz}$, 1H).

(d) 3,10,11-Tribromo-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-one

To 18.1 g (59.2 mmol) of 3-bromo-9-fluoro-10,11-dihydro-5H-dibenzo[a,d]cyclohepten-5-one and 21.4 g (120.2 mmol) of N-bromosuccineimide was added 120 ml of 1,2-dichloroethane, and the mixture was stirred at 40°C for 6 hours under mercury lamp irradiation.

After completion of the reaction, chloroform was added to the reaction mixture, and the resulting mixture was washed with a saturated aqueous sodium hydrogen carbonate solution. The organic layer was dried over anhydrous magnesium sulfate, and concentrated under reduced pressure to obtain 28.9 g of the title compound as yellowish solid. $^1\text{H-NMR}$ (δ , CDCl_3); 5.70 (d, $J=5.6\text{Hz}$, 1H), 6.14 (d, $J=5.9\text{Hz}$, 1H), 7.27-7.38 (m, 2H), 7.46-7.53 (m, 1H), 7.69 (dd, $J=8.3\text{Hz}$, 2.2Hz , 1H), 7.89-7.92 (m, 1H), 8.17 (d, $J=2.2\text{Hz}$, 1H).

(e) 3-Bromo-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-one

After dissolving 8.9 g (20 mmol) of 3,10,11-tribromo-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-one in 50 ml of

N,N-dimethylformamide under heating, the solution was cooled, and 14.4 g of sodium hydrosulfite and 18.5 g of sodium hydrogen carbonate was added to the solution and the resulting mixture was stirred at 40°C for 2 hours.

5 After completion of the reaction, warm water was added to the reaction mixture and the mixture was stirred until a temperature thereof became room temperature, and precipitated solid was collected by filtration. The filtrated product was washed with water and dried under
10 reduced pressure to obtain 5.1 g of the title compound as yellowish solid.

¹H-NMR (δ, CDCl₃); 7.08 (d, J=12.5Hz, 1H), 7.34-7.44 (m, 3H), 7.47-7.54 (m, 1H), 7.48 (dd, J=8.3Hz, 2.2Hz, 1H), 7.93-7.96 (m, 1H), 8.29 (d, J=2.2Hz, 1H).

15 (f) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-one

In 30 ml of N,N-dimethylformamide were dissolved 2.7 g (13 mmol) of 7-chloro-6-fluoro-2-vinyl quinoline and 3.0 g (10 mmol) of 3-bromo-9-fluoro-5H-dibenzo[a,d]cyclohepten-
20 5-one, and then, 1.0 g (4.6 mmol) of palladium acetate (II), 0.52 g (4.6 mmol) of triphenylphosphine and 2.2 ml (16 mmol) of triethylamine were added to the solution and the resulting mixture was stirred at 100°C for 4 hours under argon atmosphere.

25 After completion of the reaction, the reaction mixture was cooled to room temperature by allowing to stand, 50 ml of ethanol was added to the mixture and the resulting mixture was stirred, and then, precipitated solid was collected by filtration. The collected product by
30 filtration was washed with ethanol to obtain 2.2 g of the title compound as pale green yellowish brown solid.

CI-MS (m/z); 430 (M⁺+1).

EI-MS (m/z); 429 (M⁺).

¹H-NMR (δ, DMSO-d₆); 7.40 (s, 2H), 7.66-7.71 (m, 2H), 7.67
35 (d, J=16.1Hz, 1H), 7.84 (d, J=8.3Hz, 1H), 7.91-7.94 (m, 1H), 8.02 (d, J=16.1Hz, 1H), 8.02 (d, J=9.1Hz, 1H), 8.03

(d, $J=9.8\text{Hz}$, 1H), 8.20 (dd, $J=8.3\text{Hz}$, 1.7Hz , 1H), 8.23 (d, $J=7.8\text{Hz}$, 1H), 8.31 (d, $J=1.7\text{Hz}$, 1H), 8.42 (d, $J=8.8\text{Hz}$, 1H).

(g) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-ol

5 In 153 ml of tetrahydrofuran was suspended 2.20 g (5.1 mmol) of 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-one, 100 mg (4.6 mmol) of lithium borohydride was added to the suspension under ice-cooling, and the mixture was returned
10 to room temperature and stirred for 1 hour.

After completion of the reaction, the reaction mixture was ice-cooled and 50 ml of a saturated aqueous sodium hydrogen carbonate solution was poured to the mixture little by little, and after stirring at room
15 temperature for 10 minutes, the resulting mixture was extracted with ethyl acetate. The organic layer was washed with water and further washed with a saturated aqueous sodium chloride solution one time. Then, the organic layer was dried over anhydrous magnesium sulfate and concentrated
20 under reduced pressure to obtain 1.52 g of the title compound as yellowish solid.

CI-MS (m/z); 432 (M^++1).

EI-MS (m/z); 431 (M^+).

$^1\text{H-NMR}$ (δ , DMSO- d_6); 5.09 (bs, 1H), 6.36 (bs, 1H), 7.09-
25 7.16 (m, 1H), 7.24 (dd, $J=11.5\text{Hz}$, 1.5Hz , 1H), 7.32 (d, $J=11.7\text{Hz}$, 1H), 7.44-7.52 (m, 3H), 7.59 (d, $J=7.8\text{Hz}$, 1H), 7.67 (dd, $J=8.1\text{Hz}$, 1.9Hz , 1H), 7.93 (d, $J=16.1\text{Hz}$, 1H), 7.99-8.03 (m, 3H), 8.22 (d, $J=7.6\text{Hz}$, 1H), 8.38 (d, $J=8.5\text{Hz}$, 1H).

30 Reference example 9

The reaction was carried out in the same manner as in Reference example 8(a-g) to obtain the following compounds of Reference example 9(a-g).

(a) 2-[2-(4-Bromophenyl)ethenyl]-3-chlorobenzoic
35 acid

Appearance; white solid

¹H-NMR (δ, CDCl₃); 6.72 (d, J=16.6Hz, 1H), 7.31-7.44 (m, 2H), 7.51-7.61 (m, 4H), 7.66-7.70 (m, 2H), 13.25 (b, 1H).

(b) 2-[2-(4-Bromophenyl)ethyl]-3-chlorobenzoic acid

Appearance; white solid

5 CI-MS (m/z); 341 (M⁺+1).

¹H-NMR (δ, CDCl₃); 2.77-2.83 (m, 2H), 3.22-3.27 (m, 2H), 7.21 (d, J=8.3Hz, 1H), 7.36 (t, J=7.8Hz, 1H), 7.50 (d, J=8.5Hz, 1H), 7.66 (dd, J=8.1Hz, 2.2Hz, 1H), 7.75 (dd, J=7.8Hz, 1.2Hz, 1H), 13.28 (b, 1H).

10 (c) 3-Bromo-9-chloro-10,11-dihydro-5H-dibenzo[a,d]-cyclohepten-5-one

Appearance; yellowish brown solid

CI-MS (m/z); 323 (M⁺+1).

¹H-NMR (δ, CDCl₃); 3.13-3.17 (m, 2H), 3.31-3.35 (m, 2H),
15 7.12 (d, J=8.3Hz, 1H), 7.26 (t, J=8.1Hz, 1H), 7.52-7.56 (m, 2H), 7.62 (dd, J=7.8Hz, 1.5Hz, 1H), 7.97 (d, J=2.2Hz, 1H).

(d) 3,10,11-Tribromo-9-chloro-5H-dibenzo[a,d]cyclohepten-5-one

Appearance; pale yellowish solid

20 ¹H-NMR (δ, CDCl₃); 5.78 (d, J=5.9Hz, 1H), 6.35 (d, J=5.9Hz, 1H), 7.31 (d, J=8.3Hz, 1H), 7.43 (t, J=7.8Hz, 1H), 7.65 (dd, J=7.8Hz, 1.2Hz, 1H), 7.70 (dd, J=8.3Hz, 2.2Hz, 1H), 7.92 (dd, J=7.8Hz, 1.5Hz, 1H), 8.21 (d, J=2.2Hz, 1H)

(e) 3-Bromo-9-chloro-5H-dibenzo[a,d]cyclohepten-5-one
25 one

Appearance; pale yellowish solid

CI-MS (m/z); 321 (M⁺+1).

EI-MS (m/z); 320 (M⁺).

¹H-NMR (δ, CDCl₃); 7.10 (d, J=12.5Hz, 1H), 7.39-7.48 (m, 2H),
30 7.59 (d, J=12.7Hz, 1H), 7.70-7.74 (m, 2H), 7.97 (dd, J=8.1Hz, 1.2Hz, 1H), 8.18 (d, J=1.7Hz, 1H).

(f) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-9-chloro-5H-dibenzo[a,d]cyclohepten-5-one

Appearance; green yellowish solid

35 CI-MS (m/z); 448 (M⁺+1).

EI-MS (m/z); 445 (M⁺).

¹H-NMR (δ, DMSO-d₆); 7.52 (d, J=12.5Hz, 1H), 7.56 (d, J=12.7Hz, 1H), 7.62-7.69 (m, 2H), 8.82 (d, J=8.3Hz, 1H), 7.91-8.05 (m, 5H), 8.18-8.25 (m, 3H), 8.42 (d, J=9.0Hz, 1H).

5 (g) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-ol

Appearance; orange solid

CI-MS (m/z); 448 (M⁺+1).

10 ¹H-NMR (δ, DMSO-d₆); 5.05 (bs, 1H), 6.45 (bs, 1H), 7.37-7.52 (m, 7H), 7.67 (dd, J=8.1Hz, 1.7Hz, 1H), 7.69 (d, J=7.3Hz, 1H), 7.93 (d, J=16.1Hz, 1H), 7.90-8.02 (m, 2H), 8.21 (d, J=7.3Hz, 1H), 8.38 (d, J=8.8Hz, 1H)

Reference example 10

15 The reaction was carried out in the same manner as in Reference example 8(a-e) to obtain the following compounds of Reference example 10(a-e).

(a) 2-[2-(4-Bromophenyl)ethenyl]-3-iodobenzoic acid

Appearance; pale yellowish solid

CI-MS (m/z): 428 (M⁺+1).

20 EI-MS (m/z): 429 (M⁺).

¹H-NMR (δ, CDCl₃); 6.51 (d, J=16.6Hz, 1H), 7.15 (t, J=7.8Hz, 1H), 7.29 (d, J=16.4Hz, 1H), 7.53 (d, J=8.5Hz, 1H), 7.59 (d, J=8.5Hz, 1H), 7.69 (dd, J=7.8Hz, 1.2Hz, 1H), 8.07 (dd, J=7.8Hz, 1.2Hz, 1H), 13.28 (b, 1H).

25 (b) 2-[2-(4-Bromophenyl)ethyl]-3-iodobenzoic acid

Appearance; white solid

CI-MS (m/z); 431 (M⁺+1).

30 ¹H-NMR (δ, CDCl₃); 2.75-2.83 (m, 2H), 3.20-3.25 (m, 2H), 7.09 (t, J=7.8Hz, 1H), 7.24 (d, J=8.3Hz, 1H), 7.52 (d, J=8.3Hz, 1H), 7.78 (dd, J=7.8Hz, 1.5Hz, 1H), 8.06 (dd, J=7.8Hz, 1.2Hz, 1H), 13.3 (b, 1H).

(c) 3-Bromo-9-iodo-10,11-dihydro-5H-dibenzo[a,d]cyclohepten-5-one

Appearance; yellowish solid

35 CI-MS (m/z); 413 (M⁺+1).

¹H-NMR (δ, CDCl₃); 3.13-3.17 (m, 2H), 3.31-3.33 (m, 2H),

7.01 (t, J=7.8Hz, 1H), 7.11 (d, J=8.3Hz, 1H), 7.55 (dd, J=8.1Hz, 2.2Hz, 1H), 7.75 (dd, J=7.8Hz, 1.5Hz, 1H), 7.98-8.01 (m, 2H).

(d) 3,10,11-Tribromo-9-iodo-5H-dibenzo[a,d]cyclohepten-5-one

Appearance; yellowish solid

¹H-NMR (δ, CDCl₃); 5.80 (d, J=5.9Hz, 1H), 6.17 (d, J=6.1Hz, 1H), 7.14 (t, J=7.8Hz, 1H), 7.31 (d, J=8.1Hz, 1H), 7.71 (dd, J=8.3Hz, 2.2Hz, 1H), 8.10 (dd, J=7.8Hz, 1.2Hz, 1H), 7.24 (d, J=2.2Hz, 1H).

(e) 3-Bromo-9-iodo-5H-dibenzo[a,d]cyclohepten-5-one
CI-MS (m/z); 413 (M⁺+1).

¹H-NMR (δ, CDCl₃); 7.05 (d, J=12.5Hz, 1H), 7.18 (t, J=7.8 Hz, 1H), 7.35-7.42 (m, 2H), 7.72 (dd, J=8.3Hz, 2.2Hz, 1H), 8.00 (dd, J=8.1Hz, 1.0Hz, 1H), 8.13 (d, J=2.2Hz, 1H), 8.21 (dd, J=7.8Hz, 1.2Hz, 1H).

(f) 3-Bromo-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-one

In 28 ml of N,N-dimethylformamide were suspended 6.2 g (15.1 mmol) of 3-bromo-9-iodo-5H-dibenzo[a,d]cyclohepten-5-one obtained in Reference example 10(e) and 1.15 g (0.60 mmol) of copper (I) iodide, 11.61 g (60.4 mmol) of fluoro-sulfonyl(difluoro)methyl acetate was added to the suspension, and the resulting mixture was stirred at 80°C for 21 hours under argon atmosphere.

After completion of the reaction, the reaction mixture was cooled to room temperature by allowing to stand, diluted with ethyl acetate and precipitated material was filtered off, and the filtrate was successively washed with water and a saturated aqueous sodium chloride solution. The organic layer was dried over anhydrous sodium sulfate, concentrated and the residue was applied to silica gel column chromatography (eluent: hexane/ethyl acetate=9/1(volume ratio)) to obtain 4.0 g of the title compound as pale yellowish solid.

CI-MS (m/z); 355 (M⁺+1).

EI-MS (m/z); 354 (M^+).

^1H -NMR (δ , DMSO- d_6); 7.14 (d, $J=12.5\text{Hz}$, 1H), 7.39-7.43 (m, 2H), 7.61 (t, $J=8.3\text{Hz}$, 1H), 7.75 (dd, $J=8.3\text{Hz}$, 2.2Hz, 1H), 7.99 (d, $J=7.8\text{Hz}$, 1H), 8.15 (d, $J=2.2\text{Hz}$, 1H) 8.19 (d, $J=8.5\text{Hz}$, 1H).

The reaction was carried out in the same manner as in Reference example 8(f) to obtain the following compound of Reference example 10(g).

(g) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-one
Appearance; yellowish solid

CI-MS (m/z); 482 (M^++1).

EI-MS (m/z); 479 (M^+).

^1H -NMR (δ , DMSO- d_6); 7.35 (dd, $J=12.5\text{Hz}$, 1.7Hz, 1H), 7.50 (d, $J=12.7\text{Hz}$, 1H), 7.67 (d, $J=16.4\text{Hz}$, 1H), 7.79-7.87 (m, 2H), 7.97-8.04 (m, 3H), 8.16-8.24 (m, 5H), 8.42 (d, $J=8.8\text{Hz}$, 1H).

The reaction was carried out in the same manner as in Reference example 8(g) to obtain the following compound of Reference example 10(h).

(h) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-ol
Appearance; pale yellowish solid

CI-MS (m/z): 482 (M^++1).

^1H -NMR (δ , DMSO- d_6); 5.05 (bs, 1H), 6.53 (bs, 1H), 7.32-7.69 (m, 7H), 7.94 (d, $J=16.4\text{Hz}$, 1H), 7.99-8.02 (m, 3H), 8.10 (d, $J=6.8\text{Hz}$, 1H), 8.21 (d, $J=7.6\text{Hz}$, 1H), 8.79 (d, $J=8.8\text{Hz}$, 1H).

Reference example 11

The reaction was carried out in the same manner as in Reference example 8(f) to obtain the following compound of Reference example 11(a).

(a) 3-[(E)-2-(5,6,7,8-tetrahydroquinolin-2-yl)-ethenyl]-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-one
Appearance; brown solid

CI-MS (m/z); 432 (M^++1).

¹H-NMR (δ, DMSO-d₆); 1.74-1.92 (m, 4H), 2.72-2.87 (m, 4H), 7.31 (dd, J=12.7Hz, 2.2Hz, 1H), 7.36-7.53 (m, 4H), 7.66-7.83 (m, 3H), 8.07 (dd, J=8.0Hz, 1.7Hz, 1H), 8.11-8.21 (m, 3H).

5 The reaction was carried out in the same manner as in Reference example 8(g) to obtain the following compound of Reference example 11(b).

(b) 3-[(E)-2-(5,6,7,8-tetrahydroquinolin-2-yl)-ethenyl]-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-ol

10 Appearance; gray solid

CI-MS (m/z); 434 (M⁺+1).

¹H-NMR (δ, DMSO-d₆); 1.75-1.85 (m, 4H), 2.72-2.86 (m, 4H), 5.03 (bs, 1H), 6.46 (bs, 1H), 7.27 (d, J=16.1Hz, 1H), 7.33-7.46 (m, 4H), 7.54-7.67 (m, 3H), 7.94 (s, 1H) 8.08 (d, 15 J=7.3Hz, 1H).

Reference example 12

(a) 2-Chloro-6-methylbenzoic acid

To a liquid in which 42.5 g (0.43 mol) of copper (I) chloride had been suspended in 700 ml of acetonitrile was 20 added 39.9 g (0.39 mol) of t-butyl nitrite, and the mixture was warmed to 55°C. To the liquor was added 40.0 g (0.26 mol) of 2-amino-6-methylbenzoic acid by dividing into portions. After stirring at 60°C for 3 hours, the mixture was cooled to room temperature by allowing to stand.

25 After completion of the reaction, the reaction mixture was poured into a mixed solution of 1.5 liters of conc. hydrochloric acid and 1.5 liters of water, and extracted with chloroform. The organic layer was washed with a saturated aqueous ammonium chloride solution, and 30 dried over anhydrous magnesium sulfate. The concentrate concentrated under reduced pressure was recrystallized from cyclohexane to obtain 21.8 g of the title compound as beige color solid.

CI-MS (m/z); 171 (M⁺+1).

35 ¹H-NMR (δ, CDCl₃); 2.45 (s, 3H), 7.15 (t, J=4.2Hz, 1H), 7.26-7.28 (m, 2H), 9.94 (bs, 1H).

(b) Methyl 2-chloro-6-methylbenzoate

In 10 ml of N,N-dimethylformamide was dissolved 1.8 g (10.6 mmol) of 2-chloro-6-methylbenzoic acid, and then, 1.49 g (10.8 mmol) of potassium carbonate and 2.7 ml of methyl iodide were added to the solution, and the mixture was stirred at room temperature for 3 hours.

After completion of the reaction, water was added to the reaction mixture and the mixture was extracted with diethyl ether. The organic layer was successively washed with water and a saturated aqueous sodium chloride solution, dried over anhydrous magnesium sulfate, and concentrated under reduced pressure. The residue was applied to silica gel column chromatography (eluent: hexane/ethyl acetate=9/1(volume ratio)) to obtain 1.69 g of the title compound as pale yellowish liquid.

CI-MS (m/z); 171 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , CDCl_3); 2.32 (s, 3H), 3.95 (s, 3H), 7.10 (t, $J=4.3\text{Hz}$, 1H), 7.22-7.23 (m, 2H).

(c) Methyl 2-chloro-6-bromomethylbenzoate

To 20.0 g (108 mmol) of methyl 2-chloro-6-methylbenzoate and 19.3 g (108 mmol) of N-bromosuccineimide was added 100 ml of 1,2-dichloroethane, and the mixture was stirred at 50°C for 5 hours under mercury lamp irradiation.

After completion of the reaction, chloroform was added to the reaction mixture and the resulting mixture was washed with a saturated aqueous sodium hydrogen carbonate solution. The organic layer was dried over anhydrous magnesium sulfate and concentrated under reduced pressure, and hexane was added to the residue and precipitated material was removed by filtration. The filtrate was concentrated under reduced pressure to obtain 27.6 g of the title compound as pale yellowish liquid.

$^1\text{H-NMR}$ (δ , CDCl_3); 4.00 (s, 3H), 4.50 (s, 2H), 7.32-7.39 (m, 3H).

(d) 3-Chloro-2-methoxycarbonylbenzyltriphenylphosphonium bromide

To a mixture of 27.6 g of methyl 2-chloro-6-bromo-methylbenzoate and 39.0 g of triphenylphosphine was added 200 ml of acetonitrile, and the mixture was stirred at 80°C for 3 hours.

5 After completion of the reaction, the reaction mixture was concentrated under reduced pressure and diethyl ether was added to the residue. The precipitated solid was collected by filtration and dried under reduced pressure to obtain 44.6 g of the title compound as beige color solid.
10 ¹H-NMR (δ, CDCl₃); 3.66 (s, 3H), 5.62 (d, J=2.0Hz, 2H), 7.32-7.39 (m, 18H).

(e) Methyl 6-[2-(4-bromophenyl)ethenyl]-2-chlorobenzoate

To a mixture of 44.6 g (84.9 mmol) of 3-chloro-2-methoxycarbonylbenzyltriphenylphosphonium bromide and 16.7
15 g (90.4 mmol) of 4-bromobenzaldehyde was added 150 ml of acetonitrile, and then, 14.8 g (119 mmol) of 1,5-diazabicyclo[4,3,0]-5-nonene was added dropwise to the mixture. Thereafter, the mixture was refluxed for 3 hours.

20 After completion of the reaction, the reaction mixture was concentrated under reduced pressure, ethyl acetate was added to the residue and insoluble material was removed by filtration. The filtrate was successively washed with a 5% aqueous potassium hydrogen sulfate
25 solution, a saturated aqueous sodium hydrogen carbonate solution and a saturated aqueous sodium chloride solution in this order. The organic layer was dried over anhydrous magnesium sulfate and concentrated under reduced pressure, and the residue was applied to silica gel column chromatography (eluent: hexane/ethyl acetate=9/1 (volume ratio)) to
30 obtain 27.6 g of the title compound as colorless liquid.
¹H-NMR (δ, CDCl₃); 3.95, 4.08 (each s, 3H in total), 6.61 (dd, J=12.2Hz, 2.2Hz, 1H), 6.97-7.17 (m, 3H), 7.28-7.58 (m, 6H).

35 Reaction was carried out in the same manner as in Reference example 8(b) to obtain a compound of the

following Reference example 12(f).

(f) 6-[2-(4-Bromophenyl)ethyl]-2-chlorobenzoic acid

Appearance; white solid

CI-MS (m/z); 339 ($M^+ + 1$).

5 $^1\text{H-NMR}$ (δ , CDCl_3); 2.84-2.98 (m, 4H), 6.99-7.06 (m, 2H),
7.26-7.29 (m, 3H), 7.35-7.39 (m, 2H).

The reaction was carried out in the same manner as in Reference example 8(c) to obtain the following compound of Reference example 12(g).

10 (g) 3-Bromo-6-chloro-10,11-dihydro-5H-dibenzo[a,d]-
cyclohepten-5-one

Appearance; beige color solid

CI-MS (m/z): 323 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , CDCl_3); 3.06-3.11 (m, 2H), 3.16-3.20 (m, 2H),
15 7.08 (d, $J=8.3\text{Hz}$, 1H), 7.14 (dd, $J=6.8\text{Hz}$, 1.7Hz, 1H), 7.26-
7.31 (m, 2H), 7.54 (dd, $J=8.3\text{Hz}$, 2.4Hz, 1H), 8.01 (d,
 $J=2.2\text{Hz}$, 1H).

The reaction was carried out in the same manner as in Reference example 8(d) to obtain the following compound
20 of Reference example 12(h).

(h) 3,11-Dibromo-6-chloro-10-hydro-5H-dibenzo[a,d]-
cyclohepten-5-one

Appearance; yellowish brown foamy solid

$^1\text{H-NMR}$ (δ , CDCl_3); 3.37 (dd, $J=14.9\text{Hz}$, 6.6Hz, 1H), 3.74
25 (dd, $J=14.9\text{Hz}$, 2.4Hz, 1H), 5.56-5.74 (m, 1H), 7.19 (dd,
 $J=7.3\text{Hz}$, 1.2Hz, 1H), 7.33-7.43 (m, 3H), 7.68 (dd, $J=8.3\text{Hz}$,
2.2Hz, 1H), 8.98 (d, $J=2.0\text{Hz}$, 1H).

(i) 3-Bromo-6-chloro-5H-dibenzo[a,d]cyclohepten-5-
one

30 In 300 ml of ethyl acetate was dissolved 33.0 g of
3,11-dibromo-6-chloro-10-hydro-5H-dibenzo[a,d]cyclohepten-
5-one obtained in Reference example 12(h), and 150 ml of
triethylamine was added to the solution and the resulting
mixture was stirred at 80°C for 90 minutes.

35 After completion of the reaction, the reaction
mixture was cooled to room temperature by allowing to

stand, and the mixture was concentrated under reduced pressure. Ethyl acetate was added to the residue, and the mixture was successively washed with a 5% aqueous potassium hydrogen sulfate solution, a saturated aqueous sodium hydrogen carbonate solution and a saturated aqueous sodium chloride solution in this order. The organic layer was dried over anhydrous magnesium sulfate. The solvent was removed under reduced pressure, and the residue was applied to silica gel column chromatography (eluent: toluene) to obtain 8.91 g of the title compound as brown solid.

CI-MS (m/z); 321 ($M^+ + 1$).

$^1\text{H-NMR}$ (δ , CDCl_3); 6.98 (d, $J=12.0\text{Hz}$, 1H), 7.04 (d, $J=12.2\text{Hz}$, 1H), 7.35 (d, $J=8.3\text{Hz}$, 1H), 7.41-7.47 (m, 2H), 7.57 (dd, $J=7.3\text{Hz}$, 2.0Hz , 1H), 7.68 (dd, $J=8.3\text{Hz}$, 2.2Hz , 1H), 8.98 (d, $J=2.0\text{Hz}$, 1H).

The reaction was carried out in the same manner as in Reference example 8(f) to obtain a compound of the following Reference example 12(j).

(j) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-one
Appearance; yellowish brown solid

$^1\text{H-NMR}$ (δ , $\text{DMSO}-d_6$); 7.25 (d, $J=12.0\text{Hz}$, 1H), 7.30 (d, $J=12.2\text{Hz}$, 1H), 7.62-7.77 (m, 5H), 8.00-8.04 (m, 4H), 8.13 (dd, $J=8.3\text{Hz}$, 2.0Hz , 1H), 8.22 (d, $J=7.3\text{Hz}$, 1H), 8.41 (d, $J=8.8\text{Hz}$, 1H).

The reaction was carried out in the same manner as in Reference example 8(g) to obtain the following compound of Reference example 12(k).

(k) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-ol
Appearance; beige color solid

$^1\text{H-NMR}$ (δ , $\text{DMSO}-d_6$); 5.19 (d, $J=3.7\text{Hz}$, 1H), 6.49-6.53 (m, 2H), 7.10 (d, $J=11.7\text{Hz}$, 1H), 7.15 (d, $J=12.0\text{Hz}$, 1H), 7.35 (t, $J=7.8\text{Hz}$, 1H), 7.48-7.56 (m, 3H), 7.59 (d, $J=2.4\text{Hz}$, 1H), 7.70 (dd, $J=8.1\text{Hz}$, 1.5Hz , 1H), 7.85 (s, 1H), 7.95 (d, $J=16.4\text{Hz}$, 1H), 7.97-8.03 (m, 3H), 8.22 (d, $J=7.3\text{Hz}$, 1H),

8.40 (d, J=8.8Hz, 1H).

Reference example 13

(a) 3-Bromo-8-chloro-10,11-dihydro-5H-dibenzo[a,d]-cyclohepten-5-one

5 The title compound was obtained as a by-product of Reference example 12(g).

Appearance; beige color solid

CI-MS (m/z): 323 (M⁺+1).

¹H-NMR (δ, CDCl₃); 3.16 (s, 4H), 7.12 (d, J=8.3Hz, 1H),
10 7.24 (d, J=2.2Hz, 1H), 7.32 (dd, J=8.5Hz, 2.0Hz, 1H), 7.55 (dd, J=8.1Hz, 2.2Hz, 1H), 7.98 (d, J=8.3Hz, 1H), 8.15 (d, J=2.2Hz, 1H).

The reaction was carried out in the same manner as in Reference example 8(d) to obtain the following compound
15 of Reference example 13(b).

(b) 3,10,11-Tribromo-8-chloro-5H-dibenzo[a,d]cyclohepten-5-one

Appearance; yellowish brown solid

¹H-NMR (δ, CDCl₃); 5.66 (d, J=5.6Hz, 1H), 5.70 (d, J=5.6Hz,
20 1H), 7.27 (d, J=8.3Hz, 1H), 7.40 (d, J=2.2Hz, 1H), 7.47 (dd, J=8.3Hz, 2.0Hz, 1H), 7.69 (dd, J=8.1Hz, 2.2Hz, 1H), 8.06 (d, J=8.5Hz, 1H), 8.23 (d, J=2.2Hz, 1H).

The reaction was carried out in the same manner as in Reference example 8(e) to obtain the following compound
25 of Reference example 13(c).

(c) 3-Bromo-8-chloro-5H-dibenzo[a,d]cyclohepten-5-one

Appearance; beige color solid

¹H-NMR (δ, CDCl₃); 6.96 (d, J=12.1Hz, 1H), 7.03 (d,
30 J=12.1Hz, 1H), 7.41 (d, J=8.3Hz, 1H), 7.51 (dd, J=8.3Hz, 2.0Hz, 1H), 7.53 (d, J=2.0Hz, 1H), 7.74 (dd, J=8.3Hz, 2.2Hz, 1H), 8.16 (d, J=8.4Hz, 1H), 8.36 (d, J=2.2Hz, 1H).

The reaction was carried out in the same manner as in Reference example 8(f) to obtain the following compound
35 of Reference example 13(d).

(d) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-

ethenyl]-8-chloro-5H-dibenzo[a,d]cyclohepten-5-one

Appearance; green-yellowish brown solid

¹H-NMR (δ, DMSO-d₆); 7.25 (d, J=12.0Hz, 1H), 7.35 (d, J=12.2Hz, 1H), 7.66 (d, J=6.4Hz, 1H), 7.71 (dd, J=8.5Hz, 2.2Hz, 1H), 7.84 (d, J=8.3Hz, 1H), 7.93 (d, J=2.2Hz, 1H), 7.99-8.04 (m, 3H), 8.15 (d, J=8.8Hz, 1H), 8.19-8.24 (m, 2H), 8.38 (d, J=1.7Hz, 1H), 8.41 (d, J=9.0Hz, 1H).

The reaction was carried out in the same manner as in Reference example 8(g) to obtain the following compound of Reference example 13(e).

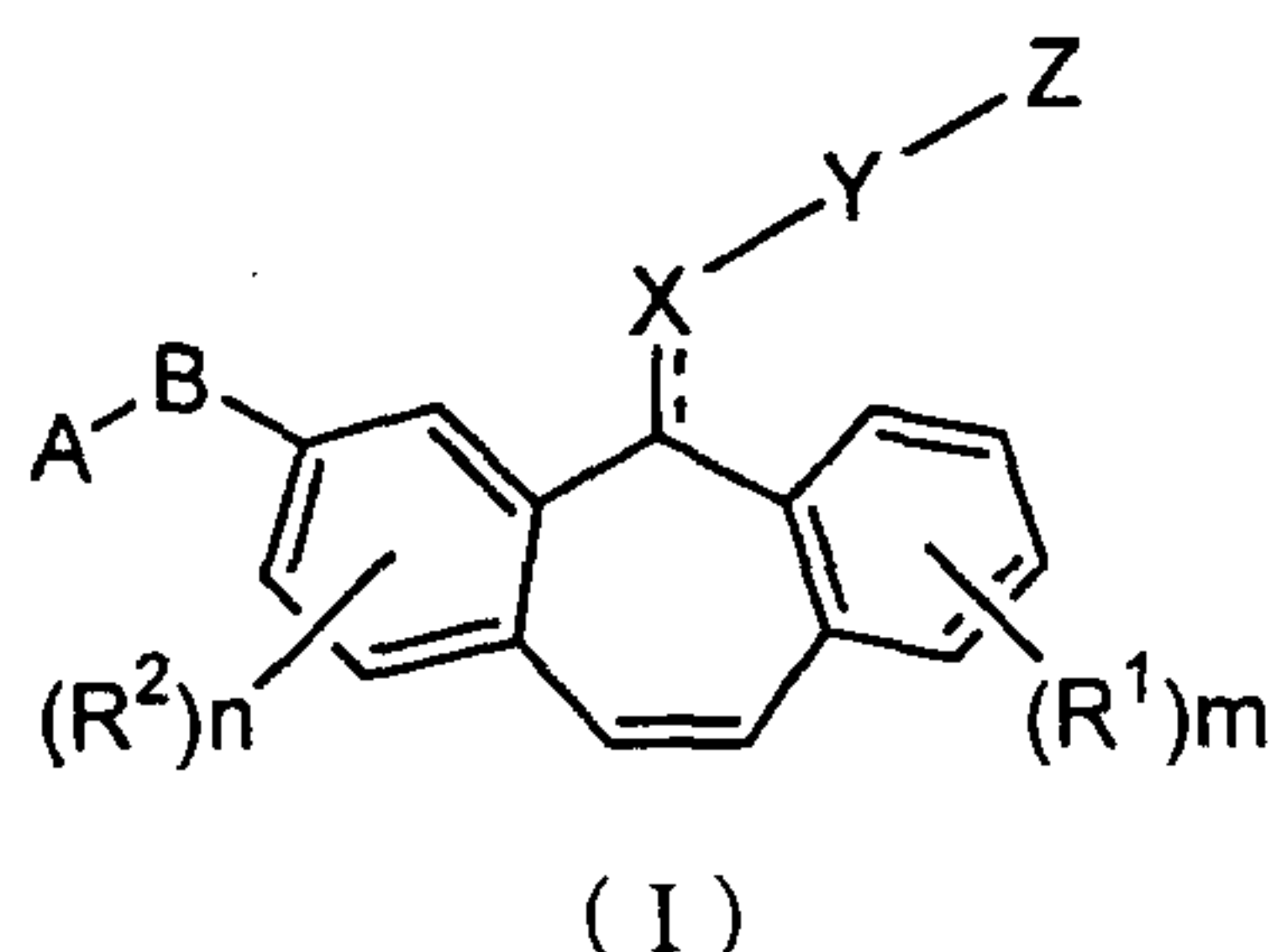
(e) 3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-ethenyl]-8-chloro-5H-dibenzo[a,d]cyclohepten-5-ol

Appearance; orange solid

¹H-NMR (δ, DMSO-d₆); 5.08 (d, J=4.2Hz, 1H), 6.32 (b, 1H), 7.15 (d, J=11.5Hz, 1H), 7.24 (d, J=11.7Hz, 1H), 7.44-7.52 (m, 4H), 7.66 (dd, J=8.1Hz, 2.0Hz, 1H), 7.75 (d, J=9.0Hz, 1H), 7.59 (d, J=16.4Hz, 1H), 7.99-8.02 (m, 3H), 8.21 (d, J=7.3Hz, 1H), 8.38 (d, J=8.8Hz, 1H).

Claims

1. A dibenzocycloheptene compound represented by the formula (I):



wherein R^1 represents a hydrogen atom, a halogen atom, a hydroxy group, a nitro group, a cyano group, a carbamoyl group, a formyl group, a carboxyl group, 1H-tetrazol-5-yl group, C_1 - C_4 alkyl group, fluoro C_1 - C_4 alkyl group, hydroxy C_1 - C_4 alkyl group, C_2 - C_4 alkenyl group, C_2 - C_4 alkynyl group, C_1 - C_4 alkoxy group, fluoro C_1 - C_4 alkoxy group, C_1 - C_4 alkylthio group, C_1 - C_4 alkylsulfinyl group or C_1 - C_4 alkylsulfonyl group,

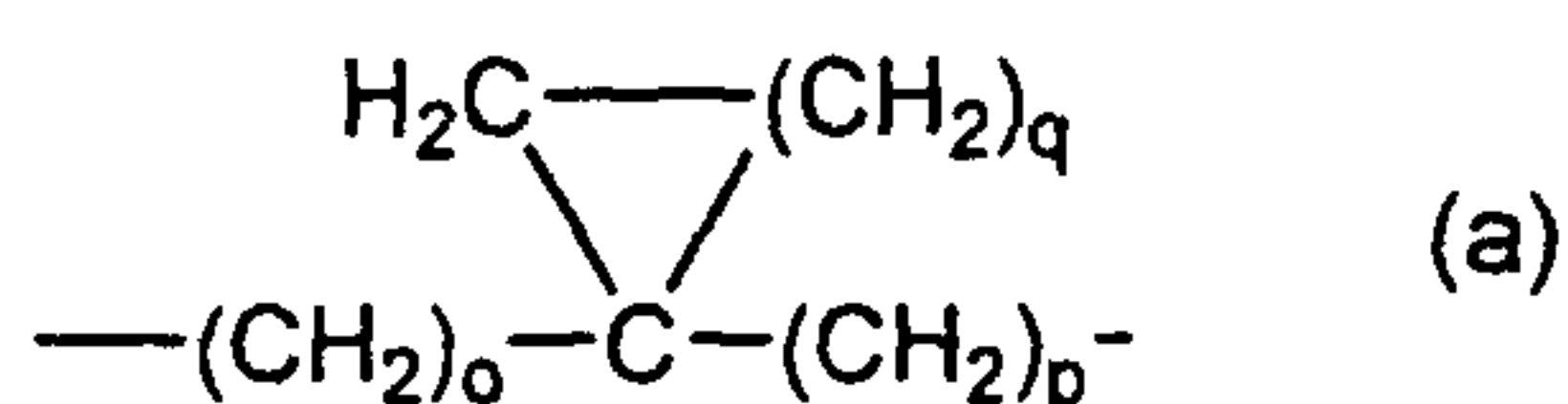
R^2 represents a hydrogen atom, a halogen atom, a nitro group, a cyano group, C_1 - C_4 alkyl group or C_1 - C_4 alkoxy group,

A represents a 5-membered or 6-membered heteroaromatic ring group having 1 to 3 hetero atom(s) selected from the group consisting of a nitrogen atom, an oxygen atom and a sulfur atom or a heteroaromatic ring-fused group in which the heteroaromatic ring group and a benzene ring are fused, and the heteroaromatic ring group or fused heteroaromatic ring group may have a halogen atom, a nitro group, a cyano group, C_1 - C_4 alkyl group, fluoro C_1 - C_4 alkyl group, C_1 - C_4 alkoxy group, fluoro C_1 - C_4 alkoxy group, C_1 - C_4 alkylthio group or C_3 - C_4 alkylene group as a substituent(s),

B represents a formula: $-\text{CH}=\text{CH}-$, a formula: $-\text{CH}_2\text{O}-$,
a formula: $-\text{CH}_2\text{CH}_2-$, a formula: $-\text{CH}_2\text{S}-$, a formula:
 $-\text{OCH}_2-$ or a formula: $-\text{SCH}_2-$,

X represents, an oxygen atom, a sulfur atom, a
methylene group or a formula: $=\text{CH}-$,

Y represents a C_1 - C_{10} alkylene group which may have a
halogen atom, C_1 - C_4 alkyl group or C_1 - C_4 alkoxy group
as a substituent(s), or a group (a) of the formula:



where o and p are each an integer of 0 to 2, and
q is an integer of 1 to 4,

Z represents a carboxyl group which may be protect-
ed; 1H-tetrazol-5-yl group; a formula: $-\text{NH}-\text{SO}_2-\text{R}^3$; or
a formula: $-\text{CO}-\text{NH}-\text{SO}_2-\text{R}^3$

where R^3 represents a C_1 - C_4 alkyl group, a fluoro-
 C_1 - C_4 alkyl group or a phenyl group which may have
a halogen atom, a C_1 - C_4 alkyl group, a fluoro- C_1 -
 C_4 alkyl group, a C_1 - C_4 alkoxy group, a fluoro- C_1 -
 C_4 alkoxy group, a nitro group or a cyano group
as a substituent(s),

m is an integer of 1 to 4, when m is 2 or more, a
plural number of R^1 may be different from each
other, n is an integer of 1 to 3, and when n is 2 or
more, a plural number of R^2 may be different from
each other,

----- represents a single bond or a double bond,
or a pharmaceutically acceptable salt thereof.

2. The dibenzocycloheptene compound or a pharmaceutically
acceptable salt thereof according to Claim 1, wherein R^1 of
the compound represented by the formula (I) is selected
from the group consisting of a hydrogen atom, a fluorine
atom, a chlorine atom, a nitro group, a cyano group, a

formyl group, a 1H-tetrazol-5-yl group, a methyl group, a difluoromethyl group, a trifluoromethyl group, a hydroxymethyl group, a 1-hydroxy-1-methylethyl group, a vinyl group, an ethynyl group, a methoxy group, a difluoromethoxy group, a trifluoromethoxy group, a methylthio group, a methylsulfinyl group and a methylsulfonyl group.

3. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to Claim 1, wherein R¹ of the compound represented by the formula (I) is selected from the group consisting of a hydrogen atom, a fluorine atom, a chlorine atom, a cyano group, a trifluoromethyl group, a hydroxymethyl group, a 1-hydroxy-1-methylethyl group, an ethynyl group, a methoxy group, a difluoromethoxy group, a trifluoromethoxy group, a methylsulfinyl group and a methylsulfonyl group.

4. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 3, wherein R² of the compound represented by the formula (I) is selected from the group consisting of a hydrogen atom, a fluorine atom, a chlorine atom, a methyl group and a methoxy group.

5. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 3, wherein R² of the compound represented by the formula (I) is a hydrogen atom.

6. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 5, wherein A of the compound represented by the formula (I) is selected from the group consisting of 2-pyridyl, 2-benzothiazolyl, quinolin-2-yl, 5,6-difluoro-2-pyridyl, 5,6-dichloro-2-pyridyl, 5,6-dimethyl-2-pyridyl, 5,6,7,8-tetrahydroquinolin-2-yl, 6-fluoro-2-benzothiazolyl, 5-

fluoro-2-benzothiazolyl, 5,6-difluoro-2-benzothiazolyl, 6-chloro-2-benzothiazolyl, 5-chloro-2-benzothiazolyl, 5,6-dichloro-2-benzothiazolyl, 5-chloro-6-fluoro-2-benzothiazolyl, 5-methyl-2-benzothiazolyl, 5-cyano-2-benzothiazolyl, 5-trifluoromethyl-2-benzothiazolyl, 5-methylthio-2-benzothiazolyl, 5-fluoroquinolin-2-yl, 6-fluoroquinolin-2-yl, 7-fluoroquinolin-2-yl, 5-chloroquinolin-2-yl, 6-chloroquinolin-2-yl, 7-chloroquinolin-2-yl, 7-methylquinolin-2-yl, 7-trifluoromethylquinolin-2-yl, 7-methoxyquinolin-2-yl, 7-difluoromethoxyquinolin-2-yl, 7-trifluoromethoxyquinolin-2-yl, 5,7-difluoroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 5,7-dichloroquinolin-2-yl, 6,7-dichloroquinolin-2-yl, 5-chloro-7-fluoroquinolin-2-yl, 6-chloro-7-fluoroquinolin-2-yl, 7-chloro-5-fluoroquinolin-2-yl, 7-chloro-6-fluoroquinolin-2-yl, 7-chloro-6-cyanoquinolin-2-yl, 7-cyano-6-fluoroquinolin-2-yl, 6-fluoro-7-trifluoromethylquinolin-2-yl and 5,6,7-trifluoroquinolin-2-yl group.

7. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 5, wherein A of the compound represented by the formula (I) is selected from the group consisting of 5,6,7,8-tetrahydroquinolin-2-yl, 7-fluoroquinolin-2-yl, 7-chloroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 6,7-dichloroquinolin-2-yl and 7-chloro-6-fluoroquinolin-2-yl group.

8. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 7, wherein B of the compound represented by the formula (I) is a formula: $-\text{CH}=\text{CH}-$, a formula: $-\text{OCH}_2-$ or a formula: $-\text{CH}_2\text{O}-$.

9. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 8, wherein X of the compound represented by the formula (I)

is an oxygen atom or a sulfur atom.

10. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 5 9, wherein Y of the compound represented by the formula (I) is selected from the group consisting of methylene, ethylene, trimethylene, difluoromethylene, 1-fluoroethylene, 2-fluoroethylene, 1,1-difluoroethylene, 2,2-difluoroethylene, ethylidene, 1-methylethylene, 2-methylethylene, 2,2-di- 10 fluorotrimethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethyltrimethylene, 2,2-dimethyltrimethylene, 3,3-dimethyltrimethylene and (a-5) group.

11. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 15 9, wherein Y of the compound represented by the formula (I) is selected from the group consisting of methylene, ethylene, trimethylene, ethylidene, 1-methylethylene, 2-methylethylene and (a-5) group.

20

12. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 11, wherein Z of the compound represented by the formula (I) is selected from the group consisting of carboxyl, 25 methanesulfonylamino, trifluoromethanesulfonylamino, methanesulfonylaminocarbonyl and trifluoromethanesulfonylaminocarbonyl group.

13. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 30 11, wherein Z of the compound represented by the formula (I) is a carboxyl group.

14. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 35 13, wherein m of the compound represented by the formula

(I) is 1 or 2.

15. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 5 14, wherein n of the compound represented by the formula (I) is 1.

16. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to Claim 1, wherein R¹ of 10 the compound represented by the formula (I) is selected from the group consisting of a hydrogen atom, a fluorine atom, a chlorine atom, a nitro group, a cyano group, a formyl group, a 1H-tetrazol-5-yl group, a methyl group, a difluoromethyl group, a trifluoromethyl group, a hydroxymethyl group, a 1-hydroxy-1-methylethyl group, a vinyl 15 group, an ethynyl group, a methoxy group, a difluoromethoxy group, a trifluoromethoxy group, a methylthio group, a methylsulfinyl group and a methylsulfonyl group, R² is selected from the group consisting of a hydrogen atom, a 20 fluorine atom, a chlorine atom, a methyl group and a methoxy group, A is selected from the group consisting of 5,6,7,8-tetrahydroquinolin-2-yl, 7-fluoroquinolin-2-yl, 7-chloroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 6,7-dichloroquinolin-2-yl and 7-chloro-6-fluoroquinolin-2-yl 25 group, B is a formula: -CH=CH-, a formula: -OCH₂- or a formula: -CH₂O-, X is an oxygen atom or a sulfur atom, Y is selected from the group consisting of methylene, ethylene, trimethylene, difluoromethylene, 1-fluoroethylene, 2-fluoroethylene, 1,1-difluoroethylene, 2,2-difluoroethylene, 30 ethylidene, 1-methylethylene, 2-methylethylene, 2,2-difluorotrimethylene, 1-methyltrimethylene, 2-methyltrimethylene, 1,1-dimethyltrimethylene, 2,2-dimethyltrimethylene, 3,3-dimethyltrimethylene and (a-5) group, Z is selected from the group consisting of carboxy, methane- 35 sulfonylamino, trifluoromethanesulfonylamino, methane-sulfonylaminocarbonyl and trifluoromethanesulfonylamino-

carbonyl group, m is 1 or 2, and n is 1.

17. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to Claim 1, wherein R¹ of
 5 the compound represented by the formula (I) is selected from the group consisting of a hydrogen atom, a fluorine atom, a chlorine atom, a cyano group, a trifluoromethyl group, a hydroxymethyl group, a 1-hydroxy-1-methylethyl group, an ethynyl group, a methoxy group, a difluoromethoxy
 10 group, a trifluoromethoxy group, a methylsulfinyl group and a methylsulfonyl group, R² is a hydrogen atom, A is selected from the group consisting of a 5,6,7,8-tetrahydroquinolin-2-yl, 7-fluoroquinolin-2-yl, 7-chloroquinolin-2-yl, 6,7-difluoroquinolin-2-yl, 6,7-dichloroquinolin-2-yl
 15 and 7-chloro-6-fluoroquinolin-2-yl group, B is a formula: -CH=CH-, a formula: -OCH₂- or a formula: -CH₂O-, X is an oxygen atom or a sulfur atom, Y is selected from the group consisting of a methylene, ethylene, trimethylene, ethyldene, 1-methylethylene, 2-methylethylene and (a-5) group, Z
 20 is a carboxy group, m is 1 or 2, and n is 1.

18. The dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to Claim 1, wherein the compound represented by the formula (I) is selected from
 25 the group consisting of:

[3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid,
 3-[[3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy]-2-methylpropionic acid,
 30 3-[[3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio]propionic acid,
 3-[[3-[(E)-2-(6,7-difluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio]-2-methylpropionic acid,
 [3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-
 35 5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid,
 3-[[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)-

ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}propionic acid,

3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}-2-methylpropionic acid,

[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thioacetic acid,

2-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thiomethyl}cyclopropane acetic acid,

[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid,

3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-9-fluoro-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-9-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid,

3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-9-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-8-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid,

2-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}propionic acid,

3-{{3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]oxy}propionic acid,

[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thioacetic acid,

3-{{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-6-chloro-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

5 3-{{[3-[(E)-2-(7-chloro-6-fluoroquinolin-2-yl)ethenyl]-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

3-{{[3-(7-chloro-6-fluoroquinolin-2-yl)methoxy-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

10 [3-[(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]oxyacetic acid,

[3-[(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thioacetic acid,

3-{{[3-[(E)-2-(7-chloroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid,

15 3-{{[3-[(E)-2-(5,6,7,8-tetrahydroquinolin-2-yl)ethenyl]-5H-dibenzo[a,d]cyclohepten-5-yl]thio}propionic acid and
or

3-{{[3-[(E)-2-(5,6,7,8-tetrahydroquinolin-2-yl)ethenyl]-9-trifluoromethyl-5H-dibenzo[a,d]cyclohepten-5-yl]thio}-
20 propionic acid.

19. A medical composition containing the dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 18 as an
25 effective ingredient.

20. The medical composition according to Claim 19, wherein the medical composition is a composition for prophylaxis or treatment of asthma, allergic diseases and/or inflammatory
30 diseases.

21. Use of the dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 18 for the preparation of a medical composition.
35 tion.

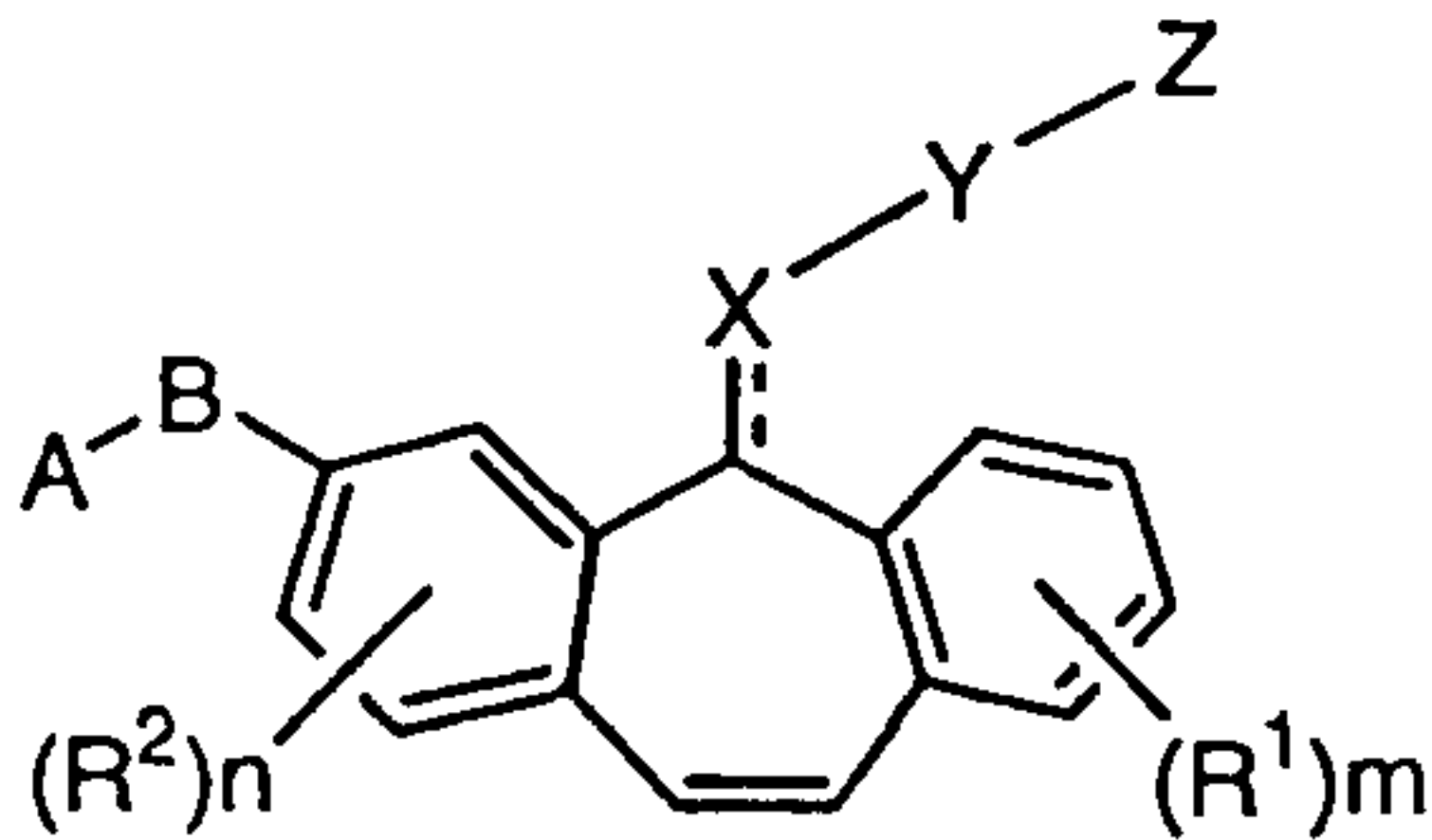
22. Use according to Claim 21, wherein the medical composition is a composition for prophylaxis or treatment of asthma, allergic diseases and/or inflammatory diseases.

5 23. A method of prophylaxis or treatment of disease which comprises administering the dibenzocycloheptene compound or a pharmaceutically acceptable salt thereof according to any one of Claims 1 to 18 to a warm-blooded animal with a pharmaceutically effective amount

10

24. The method according to Claim 23, wherein the disease is asthma, allergic diseases and/or inflammatory diseases.

15 25. The method according to Claim 23 or 24, wherein the warm-blooded animal is human.



(I)