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(54) **Novel red electroluminescent compounds and organic electroluminescent device using the same**

Neuartige rote elektrolumineszente Verbindungen und organische elektrolumineszente Vorrichtung damit

Nouveaux composés électroluminescents rouges et dispositif électroluminescent organique les utilisant

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**US-A1- 2005 227 109**

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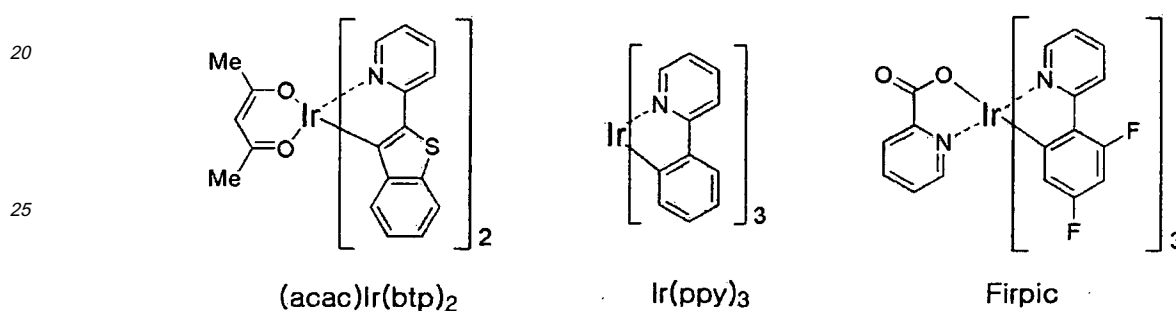
**Description**FIELD OF THE INVENTION

5 **[0001]** The present invention relates to novel organic electroluminescent compounds exhibiting high luminous efficiency, and organic electroluminescent devices and organic solar cells comprising the same.

BACKGROUND OF THE INVENTION

10 **[0002]** The most important factor to determine luminous efficiency in an OLED (organic light-emitting diode) is the type of electroluminescent material. Though fluorescent materials has been widely used as an electroluminescent material up to the present, development of phosphorescent materials is one of the best methods to improve the luminous efficiency theoretically up to four(4) times, in view of electroluminescent mechanism.

15 **[0003]** Up to now, iridium (III) complexes are widely known as phosphorescent material, including (acac)Ir(btp)<sub>2</sub>, Ir(ppy)<sub>3</sub> and Firpic, as the red, green and blue one, respectively. In particular, a lot of phosphorescent materials have been recently investigated in Japan, Europe and America.



30 **[0004]** Among conventional red phosphorescent materials, several materials have been reported to have good EL (electroluminescence) properties. However, very rare materials among them have reached the level of commercialization. As the most preferable material, an iridium complex of 1-phenyl isoquinoline may be mentioned, which is known to have excellent EL property and to exhibit color purity of dark red with high luminous efficiency. [See A. Tsuboyama et al., J. Am. Chem. Soc. 2003, 125(42), 12971-12979.]

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**[0005]** Moreover, the red materials, having no significant problem of lifetime, have tendency of easy commercialization if they have good color purity or luminous efficiency. Thus, the above-mentioned iridium complex is a material having noticeable viability of commercialization due to its excellent color purity and luminous efficiency.

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**[0006]** However, the iridium complex is still construed as a material which is merely applicable to small displays, while higher levels of EL properties than those of known materials are practically required for an OLED panel of medium to large size.

SUMMARY OF THE INVENTION

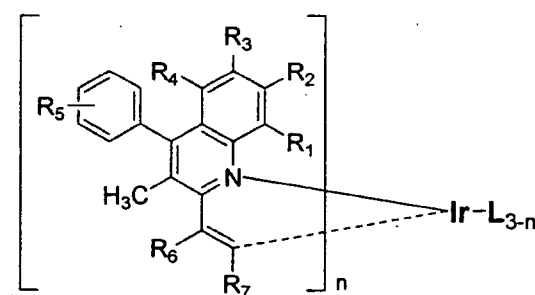
55 **[0007]** With intensive efforts to overcome the problems of conventional techniques as described above, the present inventors have researched for developing novel organic electroluminescent compounds to realize an organic EL device

having excellent luminous efficiency and surprisingly improved lifetime. Eventually, the inventors found that luminous efficiency and life property are improved when an iridium complex, which was synthesized by incorporating methyl group at 3-position of quinoline of the primary ligand compound, and a phenyl derivative at the 4-position at the same time, as a dopant of an electroluminescent device, and completed the present invention. Thus, the object of the invention is to provide novel organic electroluminescent compounds having the backbone to give more excellent properties as compared to those of conventional red phosphorescent materials. Another object of the invention is to provide novel organic electroluminescent compounds which are applicable to OLED panels of medium to large size.

[0008] Still another object of the invention is to provide organic electroluminescent devices and organic solar cells comprising the novel organic electroluminescent compounds.

[0009] The present invention relates to novel organic electroluminescent compounds and organic electroluminescent devices comprising the same. Specifically, the organic electroluminescent compounds according to the invention are characterized in that they are represented by Chemical Formula (1):

Chemical Formula 1



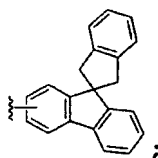
wherein, L is an organic ligand;

$R_1$  through  $R_4$  independently represent hydrogen, (C1-C60)alkyl, (C1-C60)alkoxy, (C3-C60)cycloalkyl, halogen, tri (C1-C60)alkylsilyl or tri(C6-C60)arylsilyl;

$R_5$  represents hydrogen, (C1-C60)alkyl, halogen, or (C6-C60)aryl;

$R_6$  and  $R_7$  independently represent hydrogen, (C1-C60)alkyl with or without halogen substituent(s), (C6-C60)aryl, halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino or di(C6-C60)arylamino, or  $R_6$  and  $R_7$  may be linked via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

the alkyl or aryl of  $R_6$  and  $R_7$ , or the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed therefrom by linkage via (C3-C12) alkylene or (C3-C12)alkenylene with or without a fused ring may be further substituted by one or more substituent(s) selected from (C1-C60)alkyl with or without halogen substituent(s), halogen, cyano, tri (C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino, di(C6-C60)arylamino, phenyl, naphthyl, anthryl, fluorenyl, spirobifluorenyl and



and

$n$  is an integer from 1 to 3.

#### BRIEF DESCRIPTION OF THE DRAWINGS

[0010]

Fig. 1 is a cross-sectional view of an OLED.

## DETAILED DESCRIPTION OF THE DRAWINGS

[0011] Referring now to the Drawings, Fig. 1 illustrates a cross-sectional view of an OLED comprising a Glass 1, a Transparent electrode 2, a Hole injection layer 3, a Hole transport layer 4, an Electroluminescent layer 5, an Electron transport layer 6, an Electron injection layer 7 and an Al cathode 8.

[0012] The term "alkyl" described herein and any substituents comprising "alkyl" moiety include both linear and branched species.

[0013] The term "aryl" described herein means an organic radical derived from aromatic hydrocarbon via elimination of one hydrogen atom. Each ring comprises a monocyclic or fused ring system containing from 4 to 7, preferably from 5 to 6 cyclic atoms. Specific examples include phenyl, naphthyl, biphenyl, anthryl, indenyl, fluorenyl, phenanthryl, triphenylenyl, pyrenyl, perylenyl, chrysenyl, naphthacenyl and fluoranthenyl, but they are not restricted thereto.

[0014] The term "heteroaryl" described herein means an aryl group containing from 1 to 4 heteroatom(s) selected from N, O and S as the aromatic cyclic backbone atom(s), and carbon atom(s) for remaining aromatic cyclic backbone atoms. The heteroaryl may be a 5- or 6-membered monocyclic heteroaryl or a polycyclic heteroaryl which is fused with one or more benzene ring(s), and may be partially saturated. The heteroaryl group may comprise a bivalent aryl group, of which the heteroatoms may be oxidized or quaternized to form N-oxide and quaternary salt. Specific examples include monocyclic heteroaryl groups such as furyl, thiophenyl, pyrrolyl, imidazolyl, pyrazolyl, thiazolyl, thiadiazolyl, isothiazolyl, isoxazolyl, oxazolyl, oxadiazolyl, triazinyl, tetrazinyl, triazolyl, tetrazolyl, furazanyl, pyridyl, pyrazinyl, pyrimidinyl, pyridazinyl; polycyclic heteroaryl groups such as benzofuranyl, benzothiofenyl, isobenzofuranyl, benzimidazolyl, benzothiazolyl, benzisothiazolyl, benzisoxazolyl, benzoxazolyl, isoindolyl, indolyl, indazolyl, benzothiadiazolyl, quinolyl, isoquinolyl, cinnolyl, quinazolyl, quinoxalyl, carbazolyl, phenanthridinyl and benzodioxolyl; and corresponding N-oxides (for example, pyridyl N-oxide, quinolyl N-oxide) and quaternary salts thereof; but they are not restricted thereto.

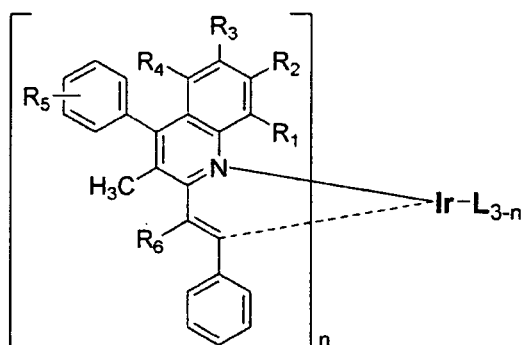
[0015] The naphthyl of Chemical Formula (1) may be 1-naphthyl or 2-naphthyl; the anthryl may be 1-anthryl, 2-anthryl or 9-anthryl; and the fluorenyl may be 1-fluorenyl, 2-fluorenyl, 3-fluorenyl, 4-fluorenyl or 9-fluorenyl.

[0016] The substituents comprising "(C1-C60)alkyl" moiety described herein may contain 1 to 60 carbon atoms, 1 to 20 carbon atoms, or 1 to 10 carbon atoms. The substituents comprising "(C6-C60)aryl" moiety may contain 6 to 60 carbon atoms, 6 to 20 carbon atoms, or 6 to 12 carbon atoms. The substituents comprising "(C3-C60)heteroaryl" moiety may contain 3 to 60 carbon atoms, 4 to 20 carbon atoms, or 4 to 12 carbon atoms. The substituents comprising "(C3-C60)cycloalkyl" moiety may contain 3 to 60 carbon atoms, 3 to 20 carbon atoms, or 3 to 7 carbon atoms. The substituents comprising "(C2-C60)alkenyl or alkynyl" moiety may contain 2 to 60 carbon atoms, 2 to 20 carbon atoms, or 2 to 10 carbon atoms.

[0017] In the compounds represented by Chemical Formula (1), the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed from R<sub>6</sub> and R<sub>7</sub> by linkage via (C3-C12) alkylene or (C3-C12)alkenylene may be benzene, naphthalene, anthracene, fluorene, indene or phenanthrene. The compound within the square bracket (□) in Chemical Formula (1) serves as a primary ligand of iridium, and L serves as a subsidiary ligand. The organic electroluminescent compounds according to the present invention also include the complex with the ratio of primary ligand: subsidiary ligand = 2:1 (n=2) and the complex with the ratio of primary ligand: subsidiary ligand = 1:2 (n=1), as well as the tris-chelated complexes without subsidiary ligand (L) (n=3).

[0018] The organic electroluminescent compounds according to the present invention can be specifically exemplified by the compounds represented by one of Chemical Formulas (2) to (7):

Chemical Formula 2

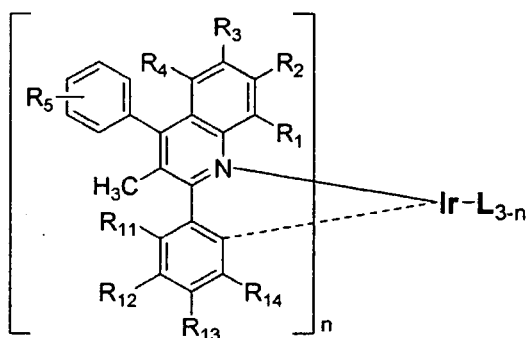


Chemical Formula 3

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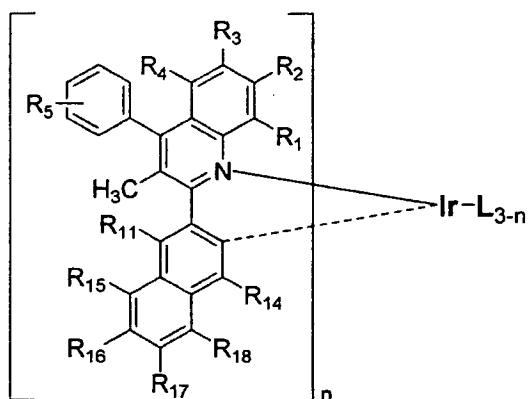


Chemical Formula 4

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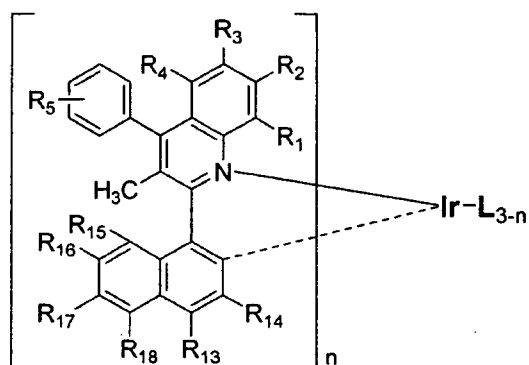
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Chemical Formula 5

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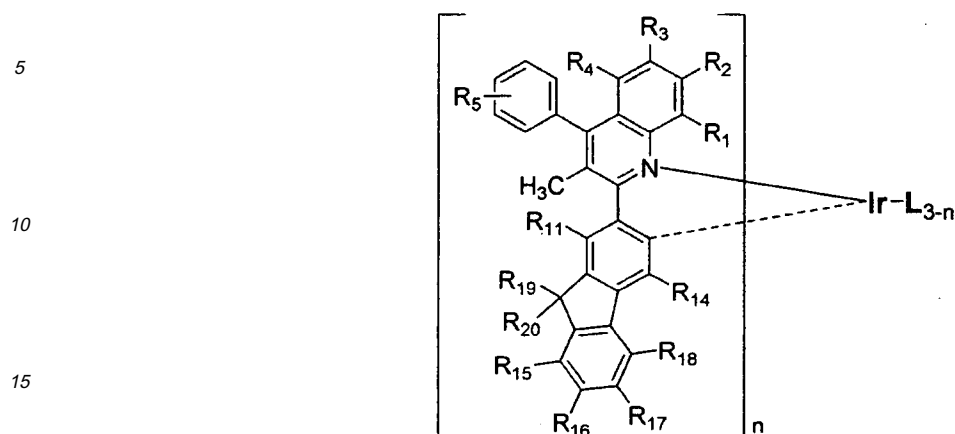
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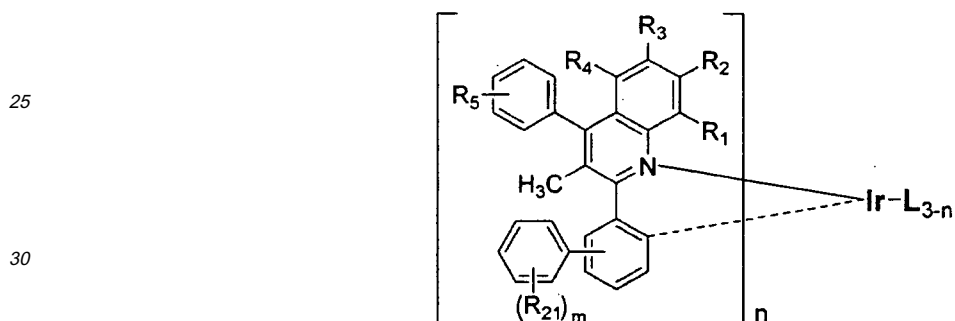


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Chemical Formula 6



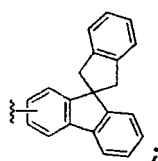
Chemical Formula 7



35 wherein, L, R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub>, R<sub>5</sub>, R<sub>6</sub> and n are defined as in Chemical Formula (1);

R<sub>11</sub> through R<sub>18</sub> independently represent hydrogen, (C1-C60)alkyl, halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino, di(C6-C60)arylamino, phenyl, naphthyl, anthryl, fluorenyl, spirobifluorenyl or

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the alkyl, phenyl, naphthyl, anthryl or fluorenyl of R<sub>11</sub> through R<sub>18</sub> may be further substituted by one or more substituent(s) selected from (C1-C60)alkyl with or without halogen substituent(s), (C1-C60)alkoxy, halogen, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino, di(C6-C60)arylamino and (C6-C60)aryl;

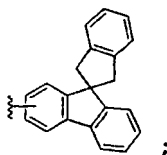
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R<sub>19</sub> and R<sub>20</sub> independently represent hydrogen, (C1-C60)alkyl or (C6-C60)aryl, or R<sub>19</sub> and R<sub>20</sub> may be linked via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring to form an alicyclic ring, a monocyclic or polycyclic aromatic ring;

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R<sub>21</sub> represents (C1-C60)alkyl, halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, phenyl, di(C1-C60)alkylamino, di(C6-C60)arylamino, naphthyl, 9,9-di(C1-C60)alkylfluorenyl or 9,9-di(C6-C60)arylfluorenyl; and m is an integer from 1 to 5.

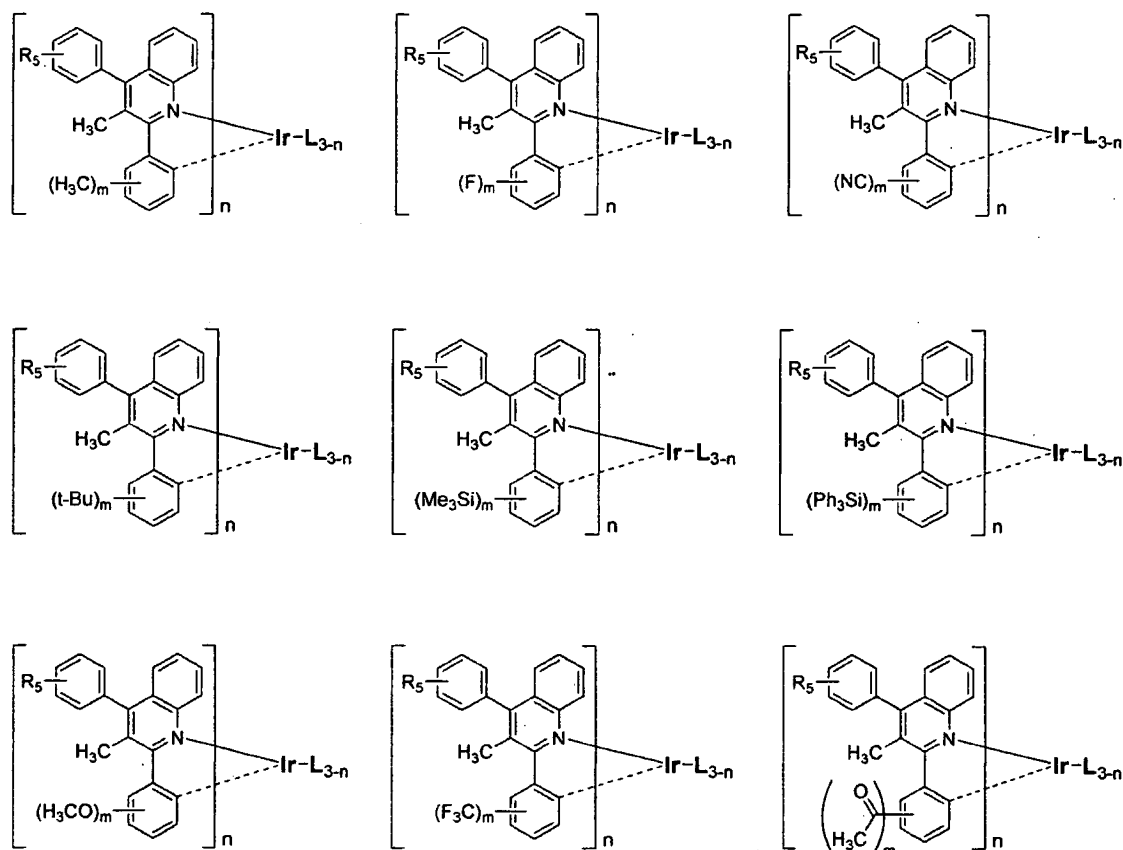
R<sub>11</sub> through R<sub>18</sub> may independently represent hydrogen, methyl, ethyl, n-propyl, i-propyl, n-butyl, i-butyl, t-butyl, n-pentyl, i-pentyl, n-hexyl, n-heptyl, n-octyl, 2-ethylhexyl, n-nonyl, trifluoromethyl, fluoro, cyano, trimethylsilyl, tripropylsilyl, tri(t-butyl)silyl, t-butyldimethylsilyl, triphenylsilyl, methoxy, ethoxy, butoxy, methylcarbonyl, ethylcarbonyl, t-butylcarbonyl, phenylcarbonyl, dimethylamino, diphenylamino, phenyl, naphthyl, anthryl, fluorenyl or



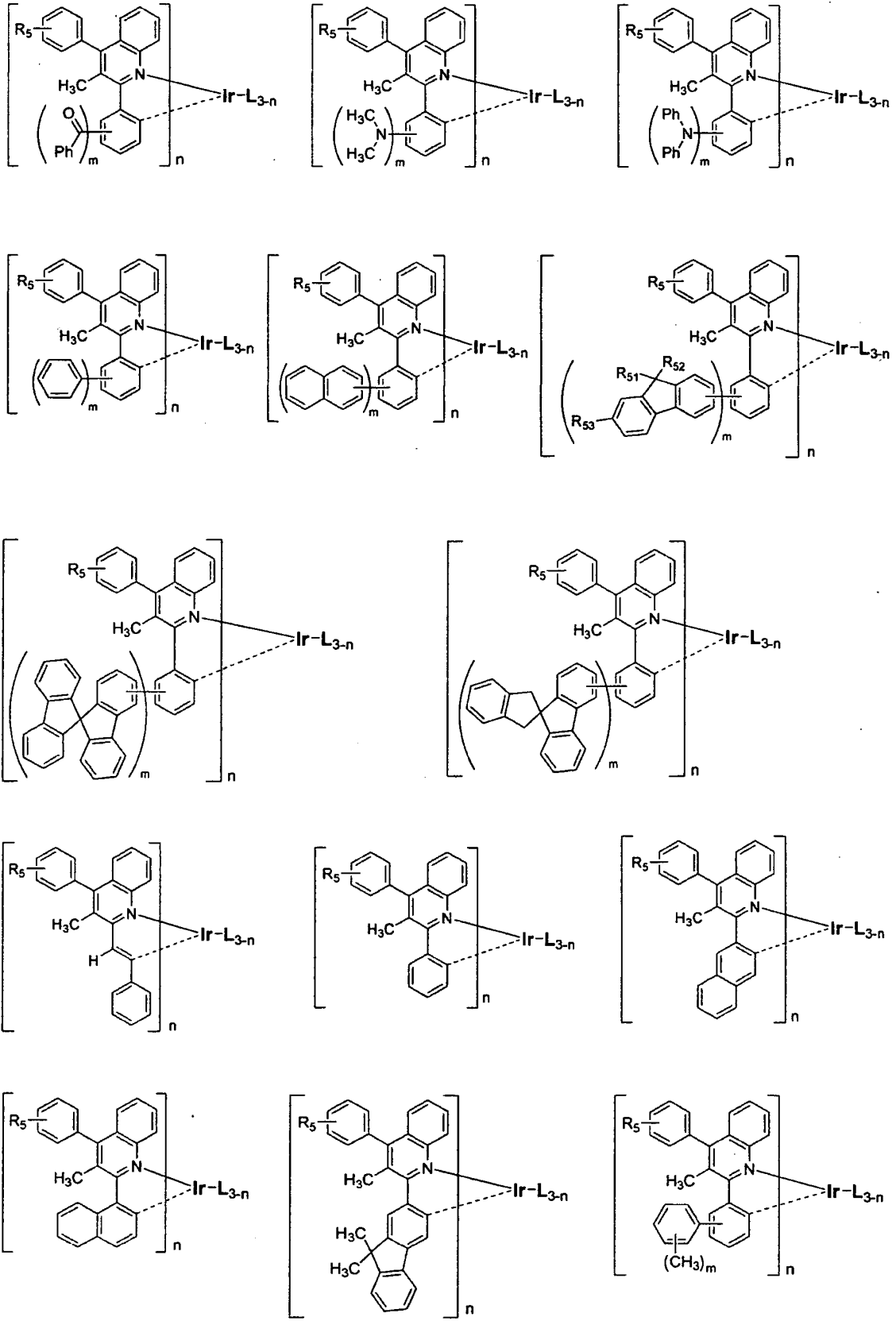
and the fluorenyl may be further substituted by methyl, ethyl, n-propyl, i-propyl, n-butyl, i-butyl, t-butyl, n-pentyl, i-pentyl, n-hexyl, n-heptyl, n-octyl, 2-ethylhexyl, n-nonyl, phenyl, naphthyl, anthryl, trimethylsilyl, tripropylsilyl, tri(t-butyl)silyl, t-butyldimethylsilyl or triphenylsilyl.

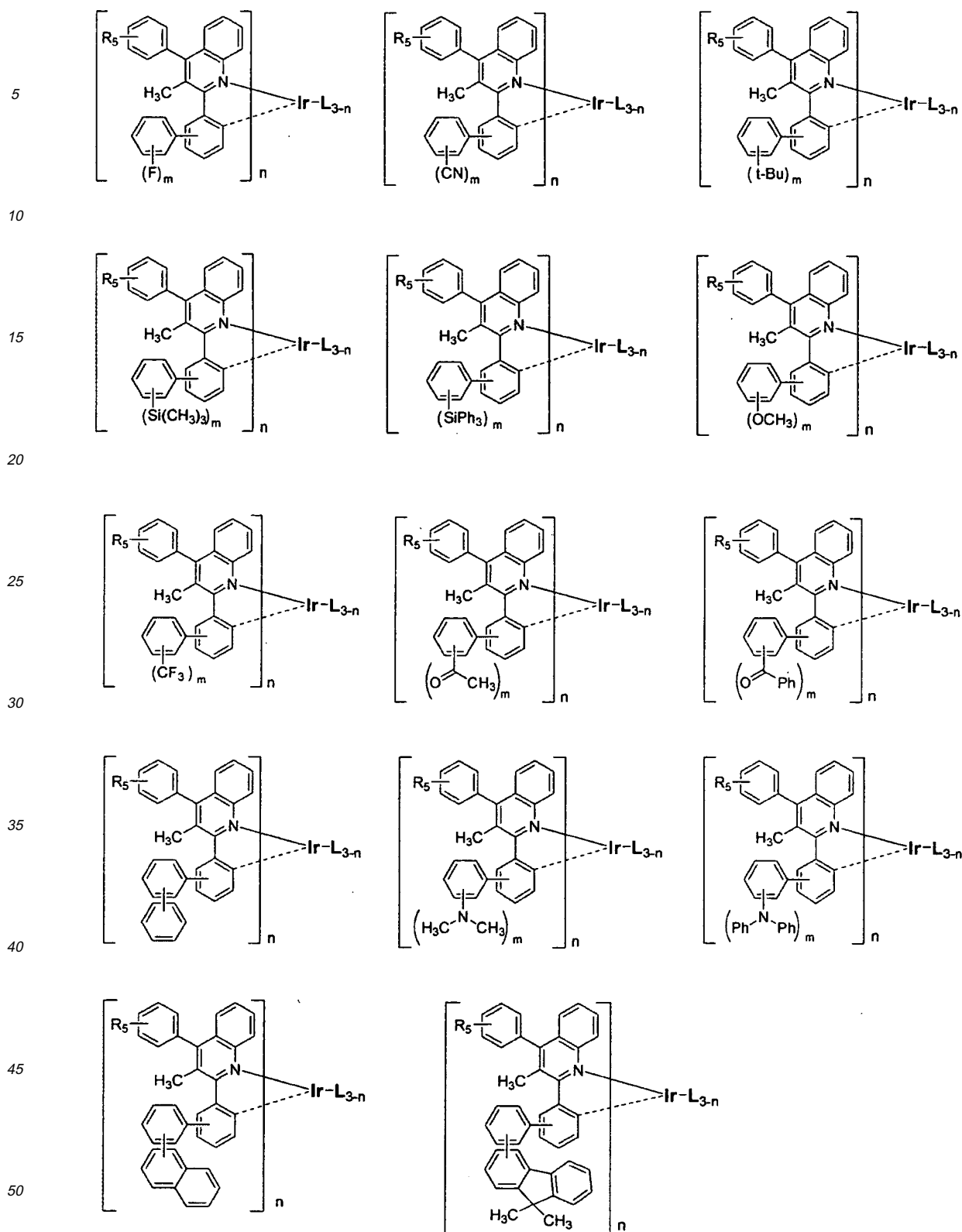
**[0019]** Preferably, R<sub>1</sub> through R<sub>4</sub> of Chemical Formula (1) independently represent hydrogen, methyl, ethyl, n-propyl, i-propyl, n-butyl, i-butyl, t-butyl, n-pentyl, i-pentyl, n-hexyl, n-heptyl, n-octyl, ethylhexyl, methoxy, ethoxy, butoxy, cyclopropyl, cyclohexyl, cycloheptyl, fluoro, trimethylsilyl, tripropylsilyl, tri(t-butyl)silyl, t-butyldimethylsilyl or triphenylsilyl; and R<sub>6</sub> represents hydrogen, methyl, ethyl, n-propyl, i-propyl, n-butyl, i-butyl, t-butyl, n-pentyl, i-pentyl, n-hexyl, n-heptyl, n-octyl, ethylhexyl, fluoro, phenyl, naphthyl, anthryl, fluorenyl or spiribifluorenyl.

**[0020]** The organic electroluminescent compounds according to the present invention can be specifically exemplified by the following compounds, but are not restricted thereto:



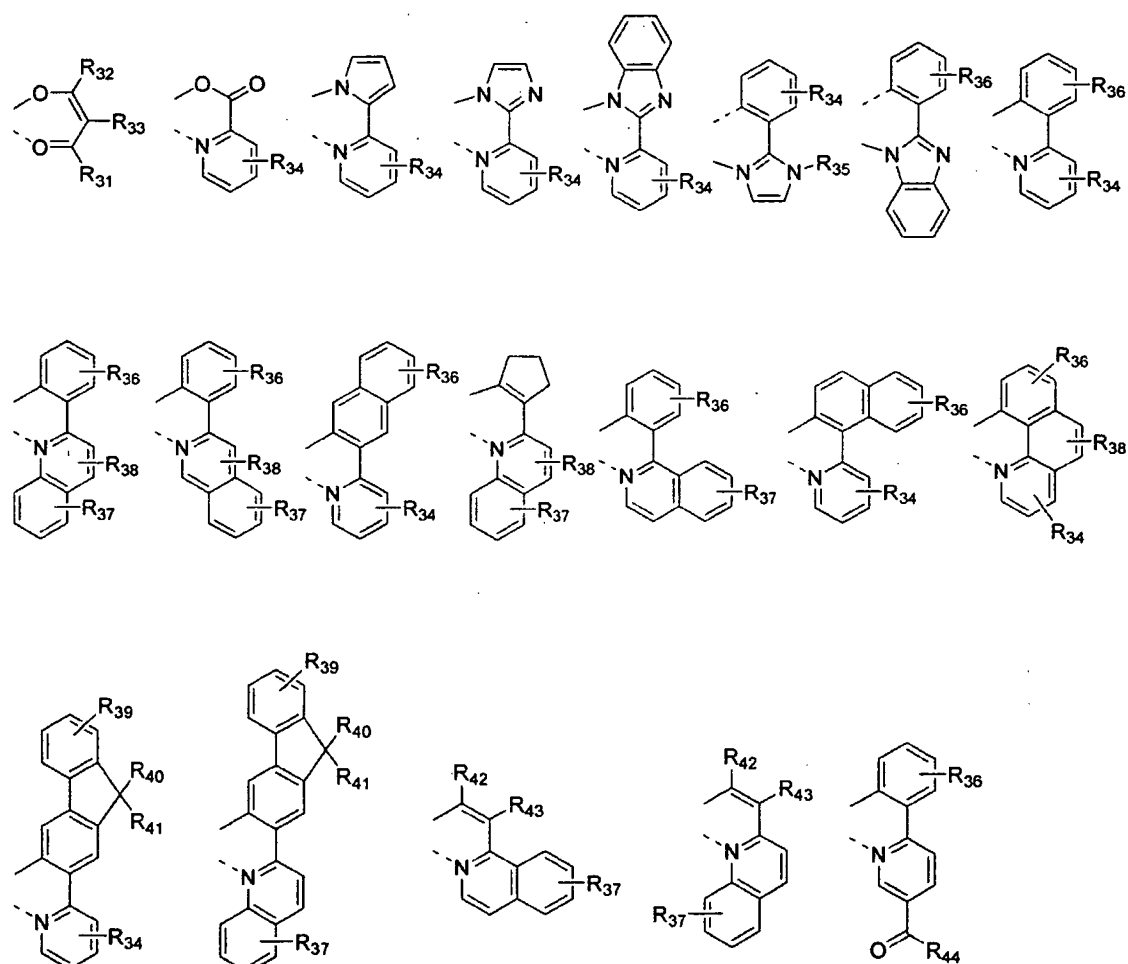
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hexyl, n-heptyl, n-octyl, ethylhexyl, phenyl or naphthyl, or  $R_{51}$  and  $R_{52}$  may be linked via (C3-C12)alkylene or (C3-C12)alkenylenes with or without a fused ring to form an alicyclic or a monocyclic or polycyclic aromatic ring;  
 $R_{53}$  represents hydrogen, methyl, ethyl, n-propyl, i-propyl, n-butyl, i-butyl, t-butyl, n-pentyl, i-pentyl, n-hexyl, n-heptyl, n-octyl, 2-ethylhexyl, trimethylsilyl, tripropylsilyl, tri(t-butyl)silyl, t-butyl dimethylsilyl, triphenylsilyl, phenyl or naphthyl;  
 m is an integer from 1 to 5; and  
 n is an integer from 1 to 3.

**[0021]** The subsidiary ligands (L) of the organic electroluminescent compounds according to the present invention include the following structures:



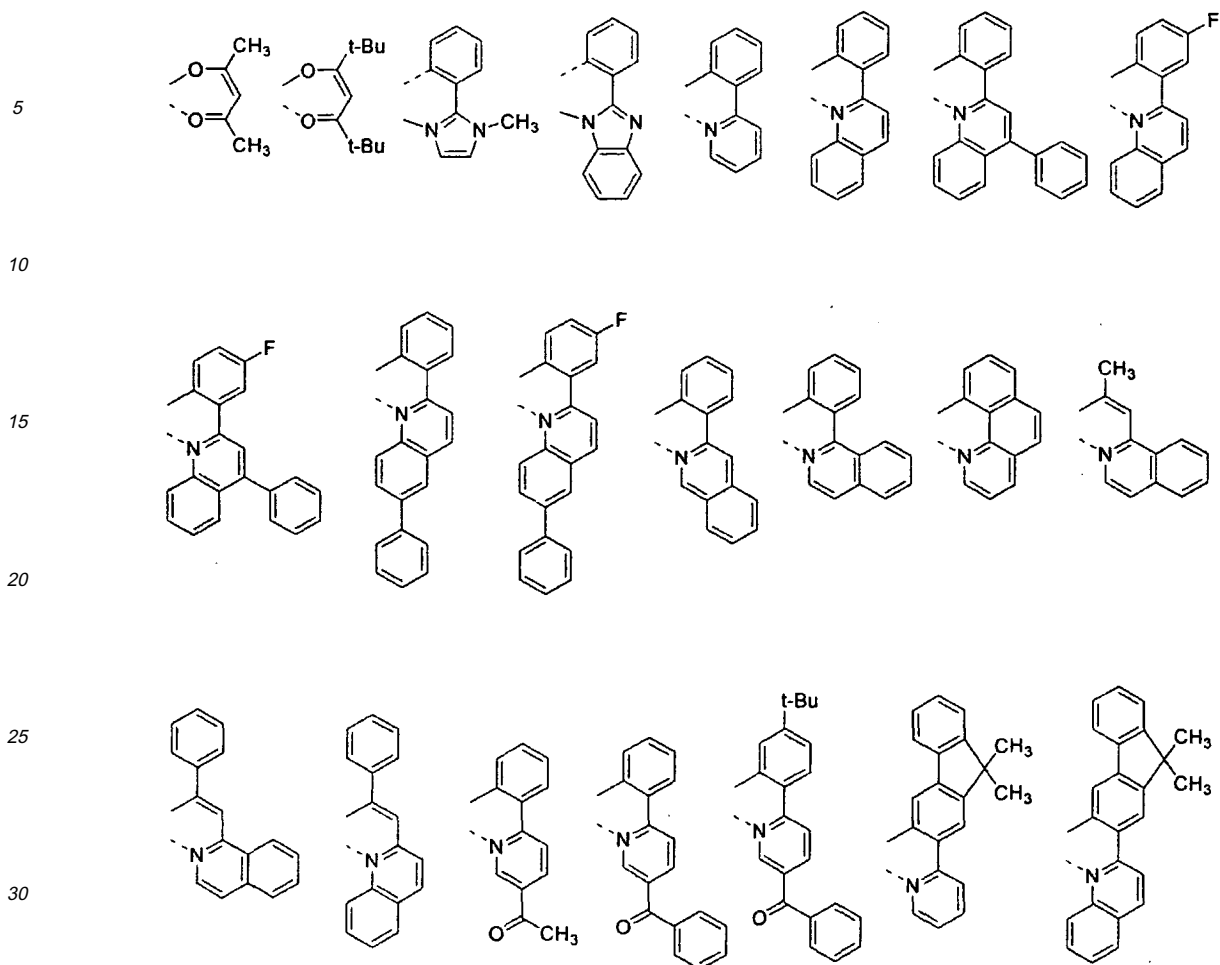
wherein,  $R_{31}$  and  $R_{32}$  independently represent hydrogen, (C1-C60)alkyl with or without halogen substituent(s), phenyl with or without (C1-C60)alkyl substituent(s), or halogen;

$R_{33}$  through  $R_{39}$  independently represent hydrogen, (C1-C60)alkyl, phenyl with or without (C1-C60)alkyl substituent(s), tri(C1-C60)alkylsilyl or halogen;

$R_{40}$  through  $R_{43}$  independently represent hydrogen, (C1-C60)alkyl, phenyl with or without (C1-C60)alkyl substituent(s); and

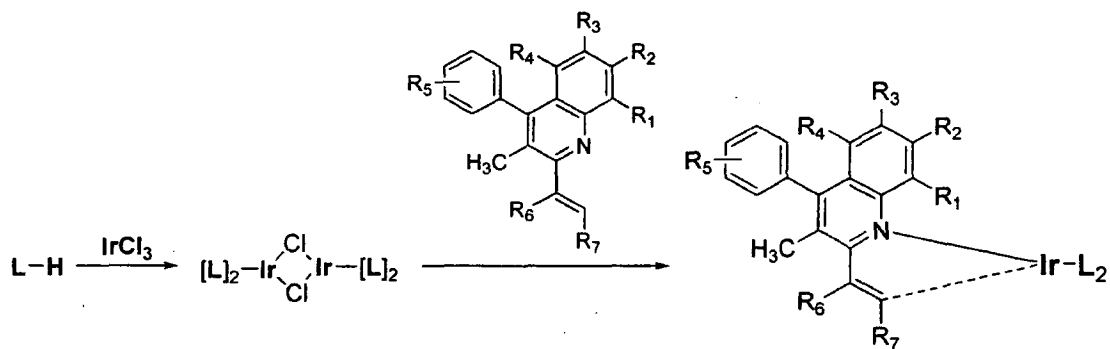
$R_{44}$  represents (C1-C60)alkyl, phenyl with or without (C1-C60)alkyl substituent(s), or halogen.

**[0022]** The subsidiary ligands (L) of the organic electroluminescent compounds according to the present invention can be exemplified by the following structures, but they are not restricted thereto:

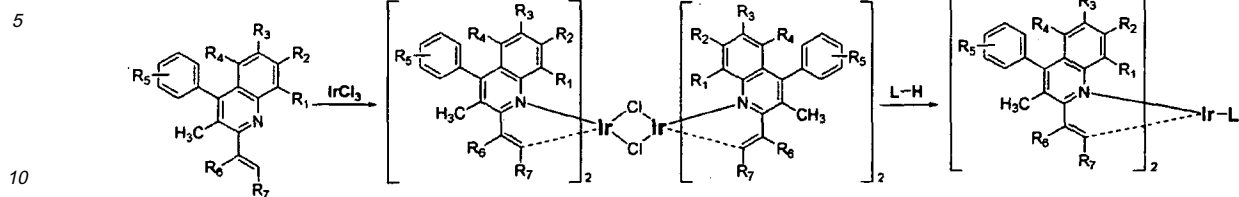


35 **[0023]** The processes for preparing the organic phosphorescent compounds according to the present invention are described by referring to Reaction Schemes (1) to (3) shown below:

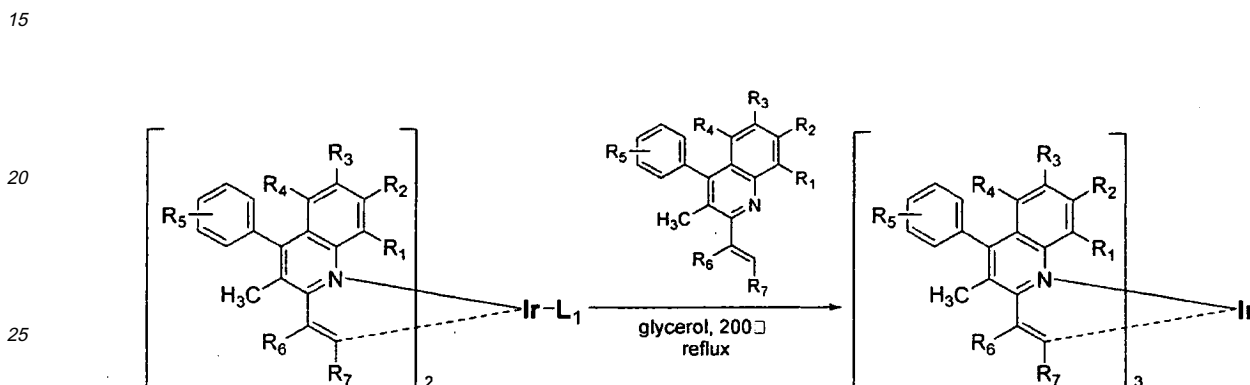
40 Reaction Scheme 1



## Reaction Scheme 2



## Reaction Scheme 3



wherein,  $R_1$ ,  $R_2$ ,  $R_3$ ,  $R_4$ ,  $R_5$ ,  $R_6$ ,  $R_7$  and  $L$  are defined as in Chemical Formula (1).

30 **[0024]** Reaction Scheme (1) provides a compound of Chemical Formula (1) with  $n=1$ , in which iridium trichloride ( $\text{IrCl}_3$ ) and subsidiary ligand compound ( $L-H$ ) are mixed in a solvent at a molar ratio of 1:2-3, and the mixture is heated under reflux before isolating diiridium dimer. In the reaction stage, preferable solvent is alcohol or a mixed solvent of alcohol/water, such as 2-ethoxyethanol, and 2-ethoxyethanol/water mixtures. The isolated diiridium dimer is then heated with a primary ligand compound in organic solvent to provide an organic phosphorescent iridium compound having the ratio of primary ligand: subsidiary ligand of 1:2 as the final product. The reaction is carried out with  $\text{AgCF}_3\text{SO}_3$ ,  $\text{Na}_2\text{CO}_3$  or  $\text{NaOH}$  being admixed with organic solvent such as 2-ethoxyethanol and 2-methoxyethylether.

35 **[0025]** Reaction Scheme (2) provides a compound of Chemical Formula (1) with  $n=2$ , in which iridium trichloride ( $\text{IrCl}_3$ ) and a primary ligand compound are mixed in a solvent at a molar ratio of 1:2-3, and the mixture is heated under reflux before isolating diiridium dimer. In the reaction stage, preferable solvent is alcohol or a mixed solvent of alcohol/water, such as 2-ethoxyethanol, and 2-ethoxyethanol/water mixtures. The isolated diiridium dimer is then heated with the subsidiary ligand compound ( $L-H$ ) in organic solvent to provide an organic phosphorescent iridium compound having the ratio of primary ligand: subsidiary ligand of 2:1 as the final product. The molar ratio of the primary ligand compound and the subsidiary ligand ( $L$ ) in the final product is determined by appropriate molar ratio of the reactant depending on the composition. The reaction may be carried out with  $\text{AgCF}_3\text{SO}_3$ ,  $\text{Na}_2\text{CO}_3$  or  $\text{NaOH}$  being admixed with organic solvent

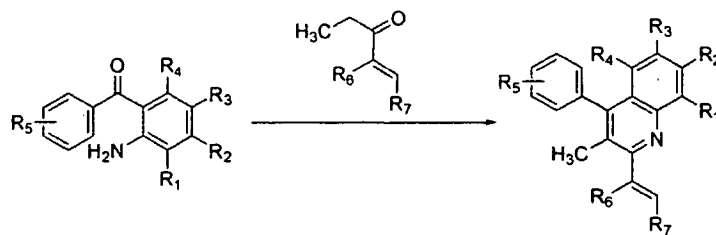
40 such as 2-ethoxyethanol, 2-methoxyethylether and 1,2-dichloroethane.

45 **[0026]** Reaction Scheme (3) provides a compound of Chemical Formula (1) with  $n=3$ , in which iridium complex prepared according to Reaction Scheme (2) and the primary ligand compound are mixed in glycerol at a molar ratio of 1:2-3, and the mixture is heated under reflux to obtain organic phosphorescent iridium complex coordinated with three primary ligands.

50 **[0027]** The compounds employed as a primary ligand in the present invention can be prepared according to Reaction Scheme (4), on the basis of conventional processes, but it is not restrictive:

## Reaction Scheme 4

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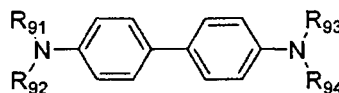
wherein,  $R_1$  through  $R_7$  are defined as in Chemical Formula (1).

**[0028]** The present invention also provides organic solar cells, which comprises one or more organic electroluminescent compound(s) represented by Chemical Formula (1).

**[0029]** The present invention also provides an organic electroluminescent device which is comprised of a first electrode; a second electrode; and at least one organic layer(s) interposed between the first electrode and the second electrode; wherein the organic layer comprises one or more compound(s) represented by Chemical Formula (1).

**[0030]** The organic electroluminescent device according to the present invention is **characterized in that** the organic layer comprises an electroluminescent region, which comprises one or more organic electroluminescent compound(s) represented by Chemical Formula (1) as electroluminescent dopant in an amount of 0.01 to 10% by weight, and one or more host(s). The host applied to the organic electroluminescent device according to the invention is not particularly restricted, but may be exemplified by 1,3,5-tricarbazolylbenzene, polyvinylcarbazole, m-biscarbazolylphenyl, 4,4'4"-tri(N-carbazolyl)triphenylamine, 1,3,5-tri(2-carbazolylphenyl)benzene, 1,3,5-tris(2-carbazolyl-5-methoxyphenyl)benzene, bis(4-carbazolylphenyl)silane or the compounds represented by one of Chemical Formulas (8) to (11).

#### Chemical Formula 8



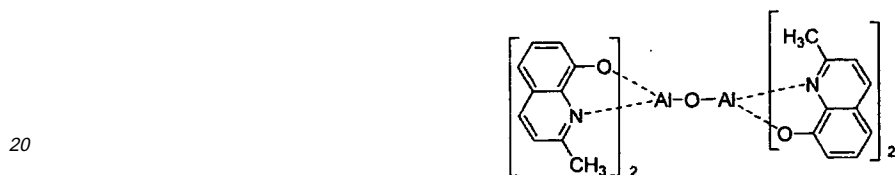
**[0031]** In Chemical Formula (8),  $R_{91}$  through  $R_{94}$  independently represent hydrogen, halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, carboxyl, nitro or hydroxyl, or each of  $R_{91}$  through  $R_{94}$  may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

the alkyl, alkenyl, alkynyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, arylsilyl, alkylsilyl, alkylamino, or arylamino of  $R_{91}$  through  $R_{94}$ , or the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed therefrom by linkage to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring may be further substituted by one or more substituent(s) selected from halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, carboxyl, nitro and hydroxyl.

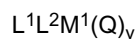
Chemical Formula 9



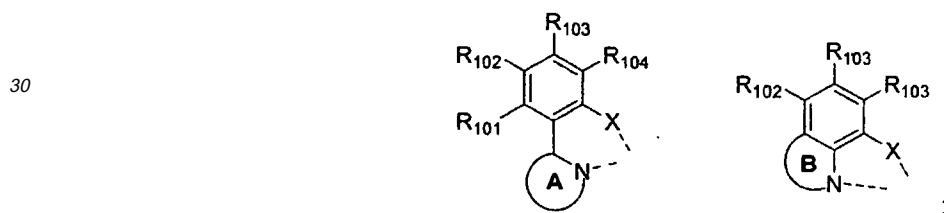
Chemical Formula 10



Chemical Formula 11



25 **[0032]** In Chemical Formula (11), the ligands,  $L^1$  and  $L^2$  are independently selected from the following structures:



$M^1$  is a bivalent or trivalent metal;

$y$  is 0 when  $M^1$  is a bivalent metal, while  $y$  is 1 when  $M^1$  is a trivalent metal;

40  $Q$  represents (C6-C60)aryloxy or tri(C6-C60)arylsilyl, and the aryloxy and triarylsilyl of  $Q$  may be further substituted by (C1-C60)alkyl or (C6-C60)aryl;

$X$  represents O, S or Se;

ring A represents oxazole, thiazole, imidazole, oxadiazole, thiadiazole, benzoxazole, benzothiazole, benzimidazole, pyridine or quinoline;

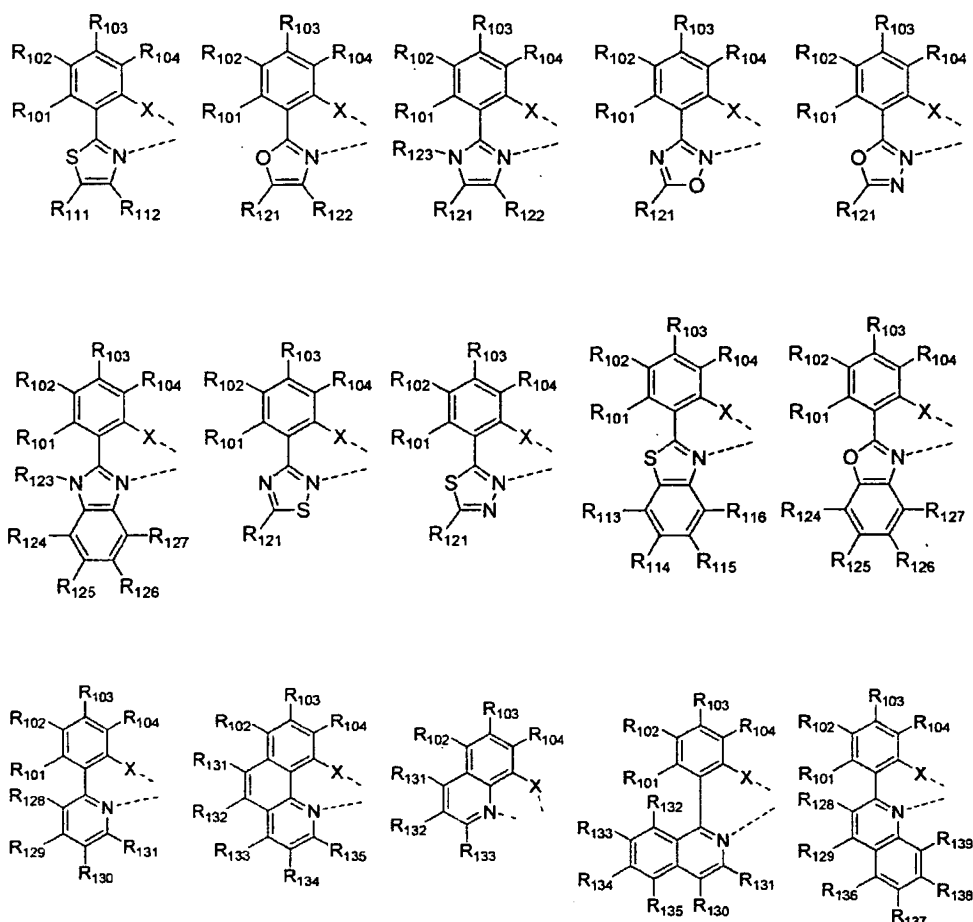
45 ring B represents pyridine or quinoline, and ring B may be further substituted by (C1-C60)alkyl, or phenyl or naphthyl with or without (C1-C60)alkyl substituent(s);

$R_{101}$  through  $R_{104}$  independently represent hydrogen, halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60) cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60) alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60) alkylcarbonyl, (C6-C60)arylcabonyl, carboxyl, nitro or hydroxyl, or each of  $R_{101}$  through  $R_{104}$  may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring; and the pyridine or quinoline may form a chemical bond with  $R_{101}$  to form a fused ring;

55 the alkyl, alkenyl, alkynyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, arylsilyl, alkylsilyl, alkylamino, or arylamino of ring A and  $R_{101}$  through  $R_{104}$ , or the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed therefrom by linkage to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring may be further substituted by one or more substituent(s) selected from halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)

heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, carboxyl, nitro and hydroxyl.

**[0033]** The ligands,  $L^1$  and  $L^2$  are independently selected from the following structures:



wherein, X represents O, S or Se;

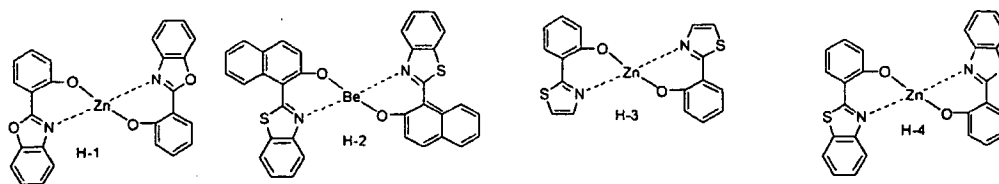
$R_{101}$  through  $R_{104}$  independently represent hydrogen, halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, carboxyl, nitro or hydroxyl, or each of  $R_{101}$  through  $R_{104}$  may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

$R_{111}$  through  $R_{116}$  and  $R_{121}$  through  $R_{139}$  independently represent hydrogen, halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, carboxyl, nitro or hydroxyl, or each of  $R_{111}$  through  $R_{116}$  and  $R_{121}$  through  $R_{139}$  may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

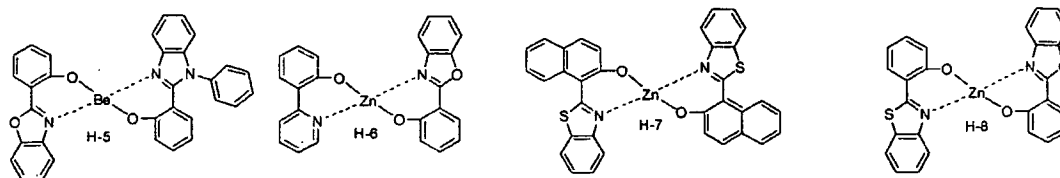


compounds with one of the following structures, but they are not restricted thereto.

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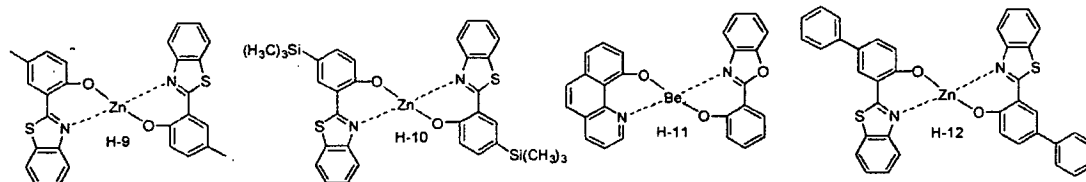


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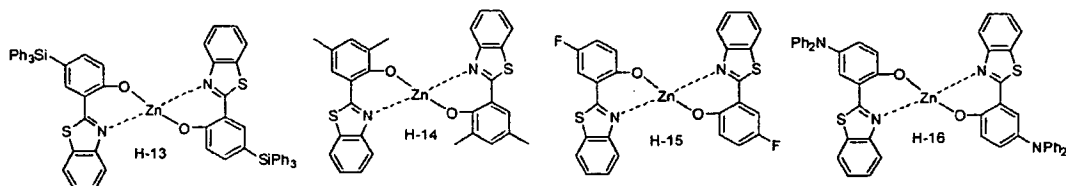
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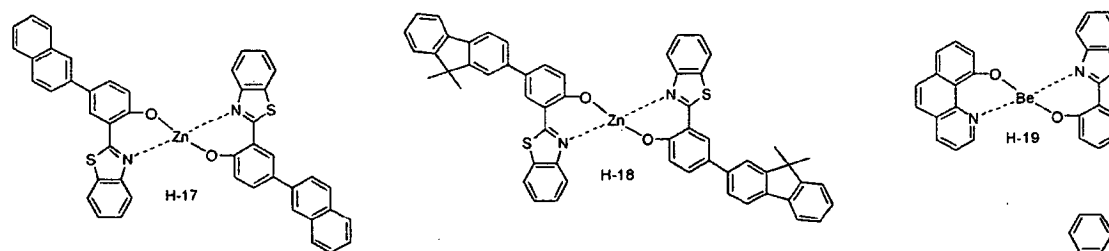
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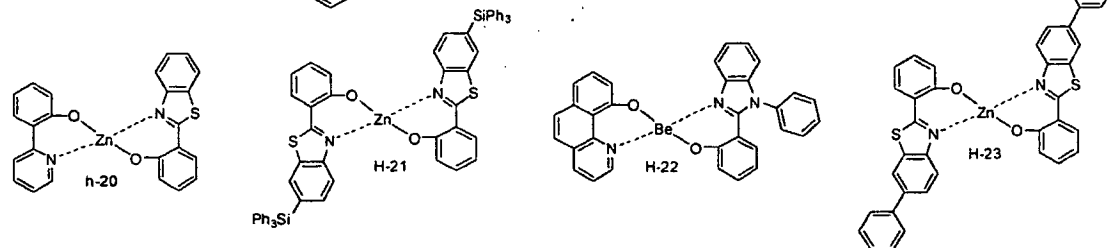


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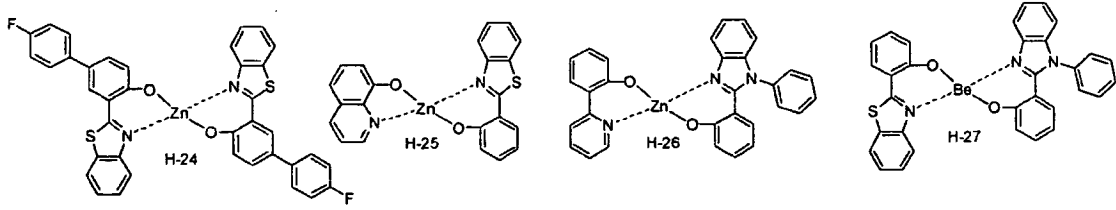
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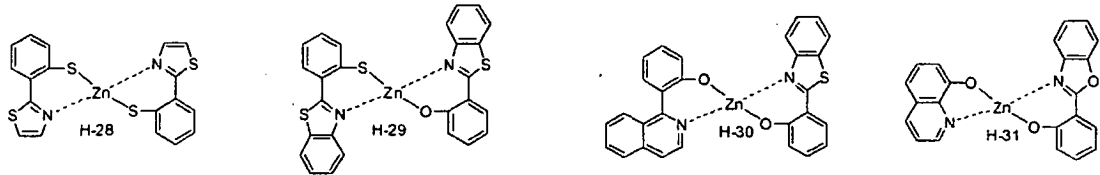
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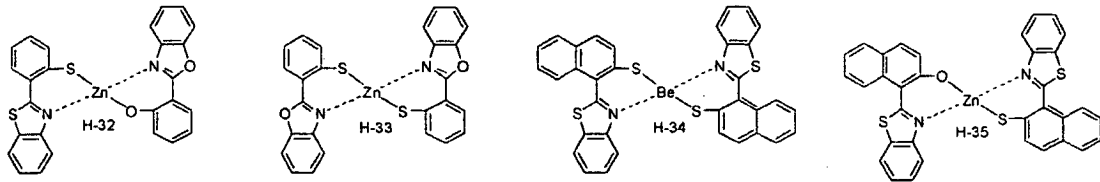
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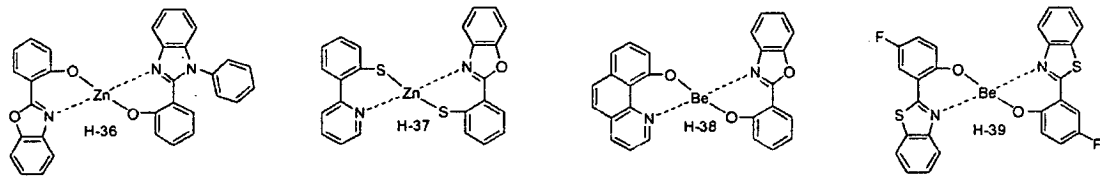


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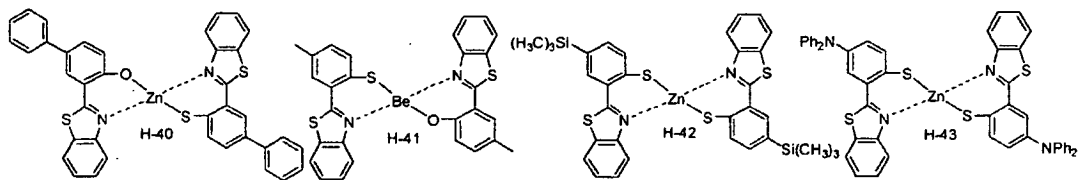


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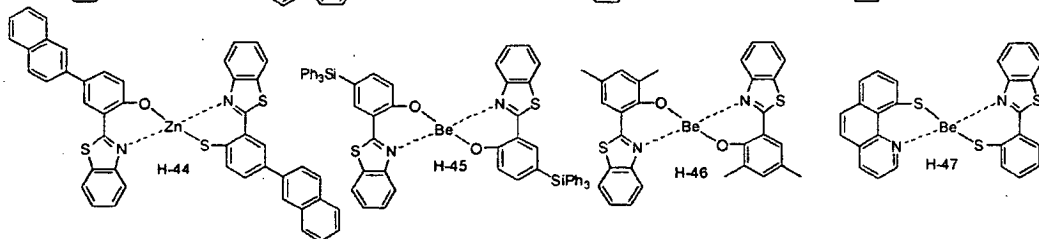


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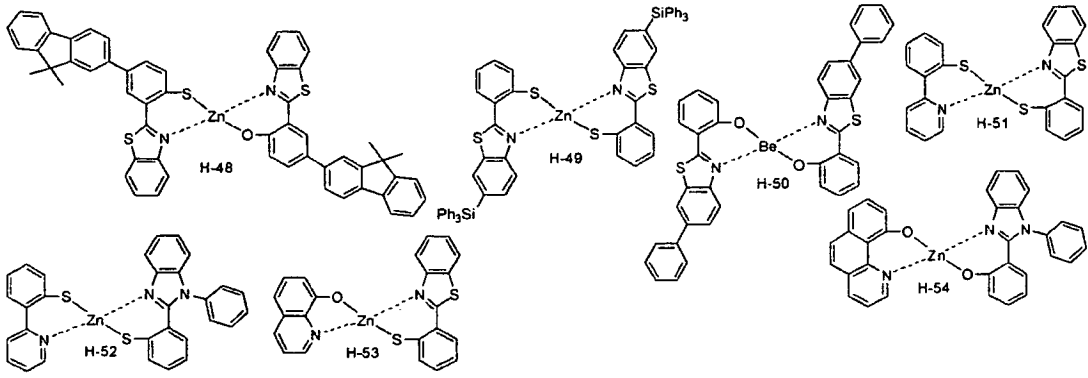
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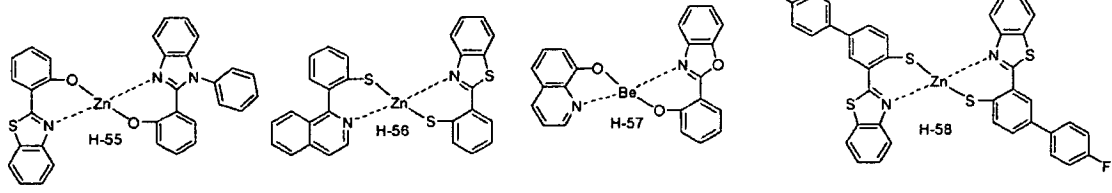
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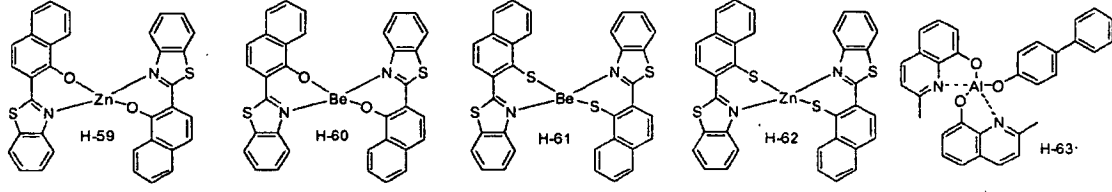
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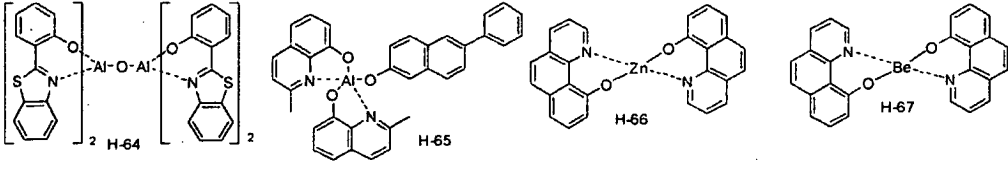
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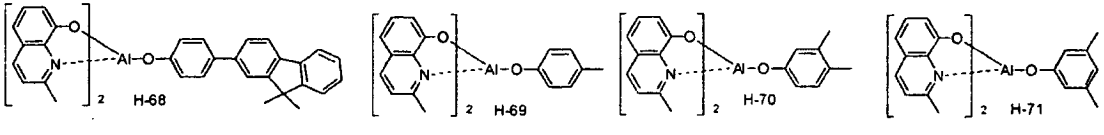


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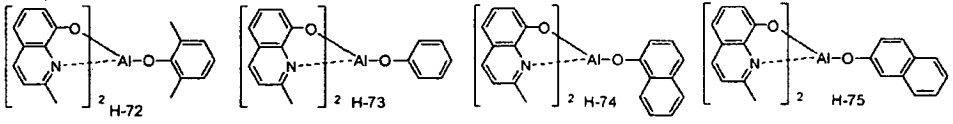
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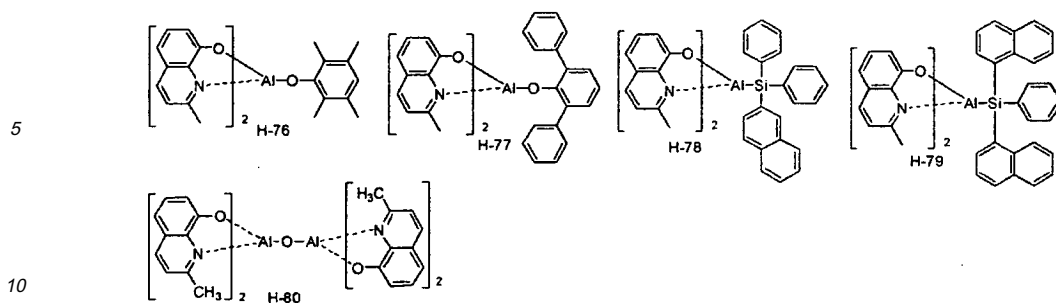
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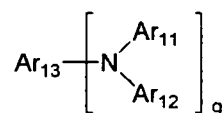
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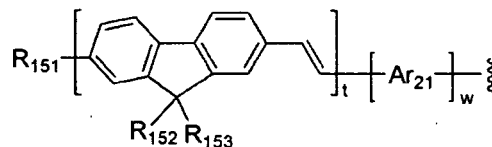
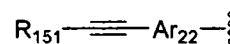
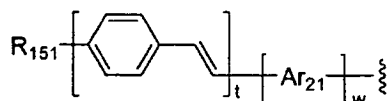
[0037] The electroluminescent layer means the layer where electroluminescence occurs, and it may be a single layer or a multi-layer consisting of two or more layers laminated. When a mixture of host-dopant is used according to the construction of the present invention, noticeable improvement in device life as well as in luminous efficiency could be confirmed.

[0038] The organic electroluminescent device according to the invention may further comprise one or more compound (s) selected from arylamine compounds and styrylarylamine compounds, as well as the organic electroluminescent compound represented by Chemical Formula (1). Examples of arylamine or styrylarylamine compounds include the compounds represented by Chemical Formula (12), but they are not restricted thereto:

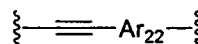
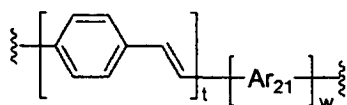
Chemical Formula 12



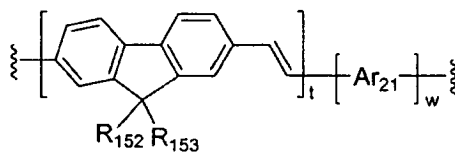
wherein,  $\text{Ar}_{11}$  and  $\text{Ar}_{12}$  independently represent (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, (C6-C60)arylamino, (C1-C60)alkylamino, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, or (C3-C60)cycloalkyl, or  $\text{Ar}_{11}$  and  $\text{Ar}_{12}$  may be linked via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring; when  $g$  is 1,  $\text{Ar}_{13}$  represents (C6-C60)aryl, (C4-C60)heteroaryl, or an aryl represented by one of the following structural formulas:



when  $g$  is 2,  $\text{Ar}_{13}$  represents (C6-C60)arylene, (C4-C60)heteroarylene, or an arylylene represented by one of the following structural formulas:



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wherein  $Ar_{21}$  and  $Ar_{22}$  independently represent (C6-C60)arylene or (C4-C60)heteroarylene;

$R_{151}$ ,  $R_{152}$  and  $R_{153}$  independently represent hydrogen, (C1-C60)alkyl or (C6-C60)aryl;

$t$  is an integer from 1 to 4,  $w$  is an integer of 0 or 1; and

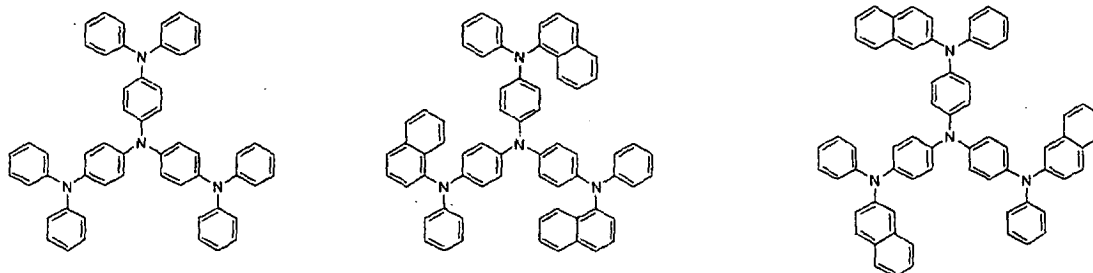
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the alkyl, aryl, heteroaryl, arylamino, alkylamino, cycloalkyl or heterocycloalkyl of  $Ar_{11}$  and  $Ar_{12}$ , or the aryl, heteroaryl, arylene or heteroarylene of  $Ar_{13}$ , or the arylene or heteroarylene of  $Ar_{21}$  and  $Ar_{22}$ , or the alkyl or aryl of  $R_{151}$  through  $R_{153}$  may be further substituted by one or more substituent(s) selected from a group consisting of halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C6-C60)aryloxy, (C1-C60)alkyloxy, (C6-C60)arylthio, (C1-C60)alkylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, carboxyl, nitro and hydroxyl.

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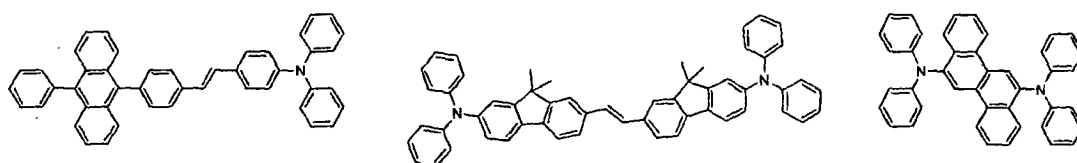
**[0039]** The arylamine compounds and styrylarylamino compounds may be more specifically exemplified by the following compounds, but are not restricted thereto.

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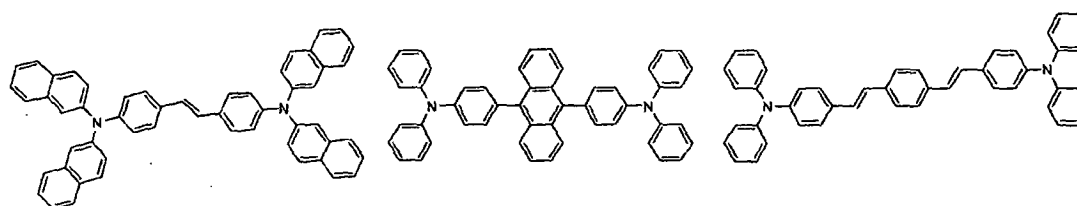
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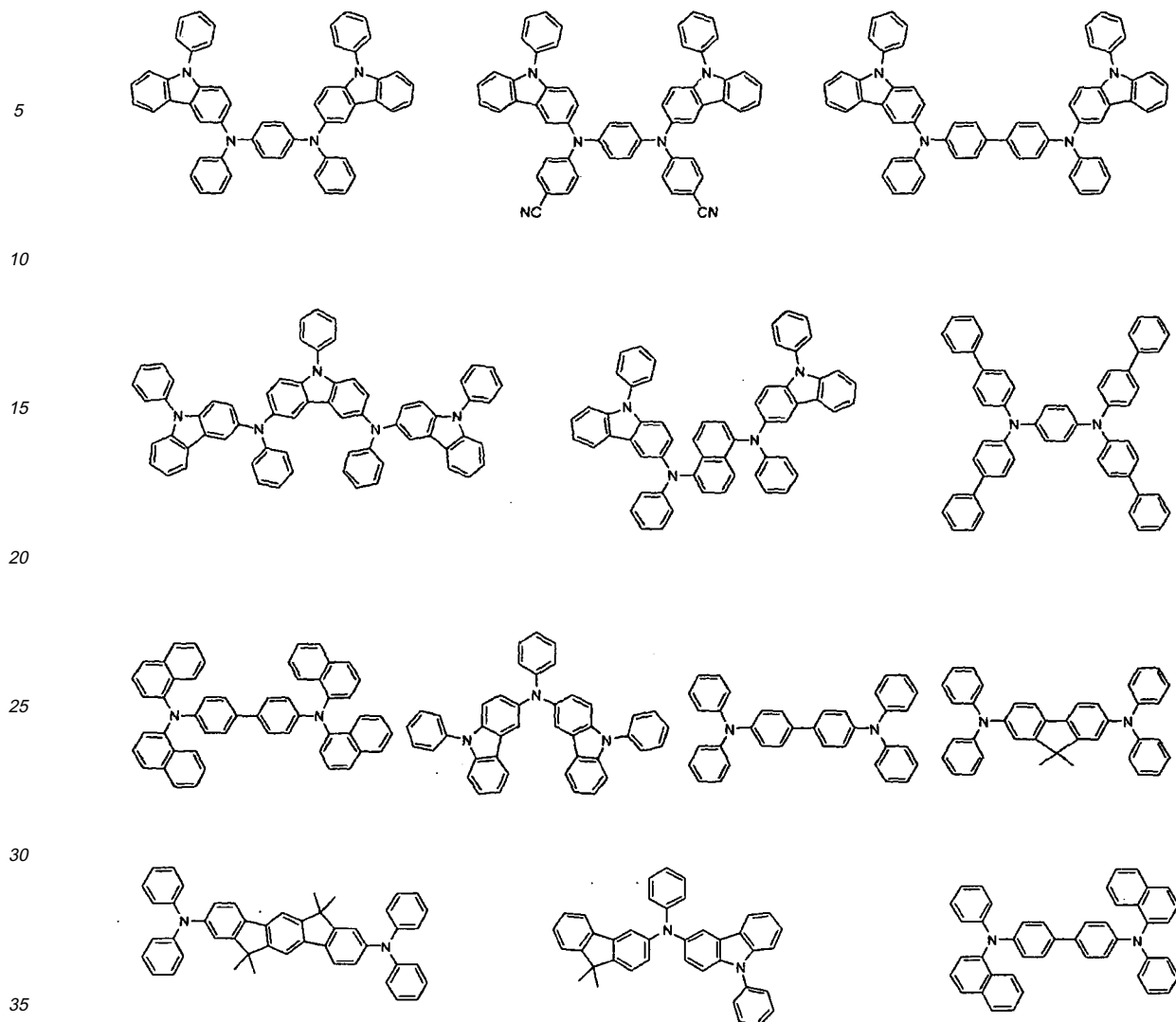
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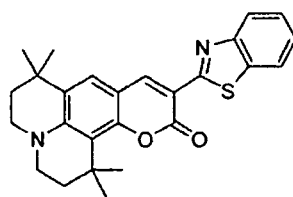


**[0040]** In an organic electroluminescent device according to the present invention, the organic layer may further comprise one or more metal(s) selected from a group consisting of organic metals of Group 1, Group 2, 4<sup>th</sup> period and 5<sup>th</sup> period transition metals, lanthanide metals and d-transition elements, as well as the organic electroluminescent compound represented by Chemical Formula (1). The organic layer may comprise a charge generating layer in addition to the electroluminescent layer.

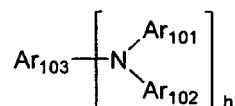
**[0041]** The present invention can realize an electroluminescent device having a pixel structure of independent light-emitting mode, which comprises an organic electroluminescent device containing the compound of Chemical Formula (1) as a sub-pixel, and one or more sub-pixel(s) comprising one or more compound(s) selected from a group consisting of arylamine compounds and styrylamine compounds, patterned in parallel at the same time.

**[0042]** Further, the organic electroluminescent device is an organic display which comprises one or more compound (s) selected from compounds having electroluminescent peak of wavelength of blue or green, at the same time. The compounds having electroluminescent peak of wavelength of blue or green may be exemplified by the compounds represented by one of Chemical Formulas (13) to (17), but they are not restricted thereto.

Chemical Formula 13

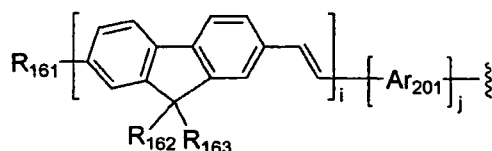
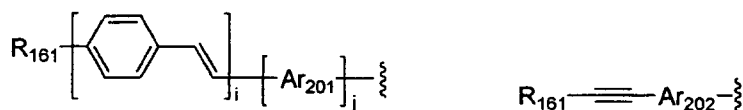


Chemical Formula 14

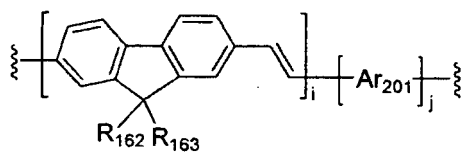
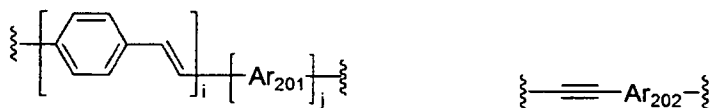


**[0043]** In Chemical Formula (14),  $\text{Ar}_{101}$  and  $\text{Ar}_{102}$  independently represent (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, (C6-C60)arylamino, (C1-C60)alkylamino, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, or (C3-C60)cycloalkyl, or  $\text{Ar}_{101}$  and  $\text{Ar}_{102}$  may be linked via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

when  $h$  is 1,  $\text{Ar}_{103}$  represents (C6-C60)aryl, (C4-C60)heteroaryl, or an aryl represented by one of the following structural formulas:



when  $h$  is 2,  $\text{Ar}_{103}$  represents (C6-C60)arylene, (C4-C60)heteroarylene, or an arylylene represented by one of the following structural formulas:



wherein  $\text{Ar}_{201}$  and  $\text{Ar}_{202}$  independently represent (C6-C60)arylene or (C4-C60)heteroarylene;

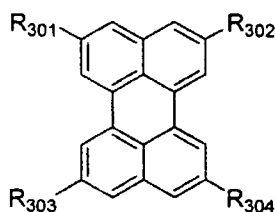
EP 2 066 150 B1

R<sub>161</sub>, R<sub>162</sub> and R<sub>163</sub> independently represent hydrogen, (C1-C60)alkyl or (C6-C60)aryl;

i is an integer from 1 to 4, j is an integer of 0 or 1; and

the alkyl, aryl, heteroaryl, arylamino, alkylamino, cycloalkyl or heterocycloalkyl of Ar<sub>101</sub> and Ar<sub>102</sub>, or the aryl, heteroaryl, arylene or heteroarylene of Ar<sub>103</sub>, or the arylene or heteroarylene of Ar<sub>201</sub>, and Ar<sub>202</sub>, or the alkyl or aryl of R<sub>161</sub> through R<sub>163</sub> may be further substituted by one or more substituent(s) selected from a group consisting of halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C6-C60)aryloxy, (C1-C60)alkyloxy, (C6-C60)arylthio, (C1-C60)alkylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcabonyl, carboxyl, nitro and hydroxyl.

Chemical Formula 15



**[0044]** In Chemical Formula (15), R<sub>301</sub> through R<sub>304</sub> independently represent hydrogen, halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcabonyl, carboxyl, nitro or hydroxyl, or each of R<sub>301</sub> through R<sub>304</sub> may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

the alkyl, alkenyl, alkynyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, arylsilyl, alkylsilyl, alkylamino or arylamino of R<sub>301</sub> through R<sub>304</sub>, or the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed therefrom by linkage to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring may be further substituted by one or more substituent(s) selected from halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcabonyl, carboxyl, nitro and hydroxyl.

Chemical Formula 16  $(Ar_{301})_p-L_{11}-(Ar_{302})_q$

Chemical Formula 17  $(Ar_{303})_r-L_{12}-(Ar_{304})_s$

**[0045]** In Chemical Formulas (16) and (17),

L<sub>11</sub> represents (C6-C60)arylene or (C4-C60)heteroarylene;

L<sub>12</sub> represents anthracenylene;

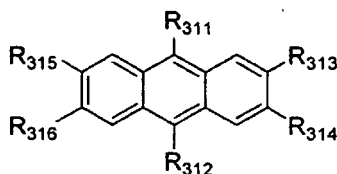
Ar<sub>301</sub> through Ar<sub>304</sub> are independently selected from hydrogen, (C1-C60)alkyl, (C1-C60)alkoxy, halogen, (C4-C60)heteroaryl, (C5-C60)cycloalkyl and (C6-C60)aryl, and the cycloalkyl, aryl or heteroaryl of Ar<sub>301</sub> through Ar<sub>304</sub> may be further substituted by one or more substituent(s) selected from a group consisting of (C6-C60)aryl or (C4-C60)heteroaryl with or without at least one substituent(s) selected from a group consisting of (C1-C60)alkyl with or without halogen substituent(s), (C1-C60)alkoxy, (C3-C60)cycloalkyl, halogen, cyano, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl and tri(C6-C60)arylsilyl; (C1-C60)alkyl with or without halogen substituent(s), (C1-C60)alkoxy,

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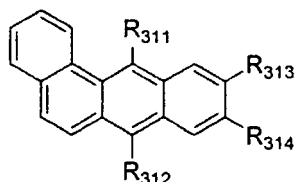
(C3-C60)cycloalkyl, halogen, cyano, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl and tri(C6-C60)arylsilyl;  
and  
p, q, r and s independently represent an integer from 0 to 4.

5 **[0046]** The compounds represented by Chemical Formula (16) or (17) may be exemplified by the derivatives represented by one of Chemical Formulas (18) to (21).

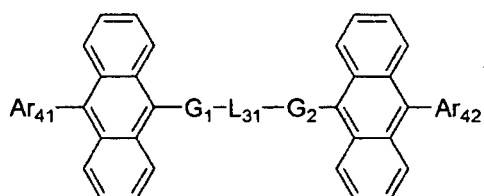
10 Chemical Formula 18



20 Chemical Formula 19



30 Chemical Formula 20



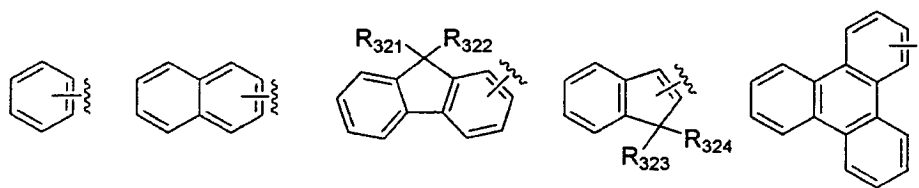
40 **[0047]** In Chemical Formulas (18) to (20), R<sub>311</sub> and R<sub>312</sub> independently represent (C6-C60)aryl, (C4-C60)heteroaryl or a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, or (C3-C60) cycloalkyl, and the aryl or heteroaryl of R<sub>311</sub> and R<sub>312</sub> may be further substituted by one or more substituent(s) selected from a group consisting of (C1-C60)alkyl, halo(C1-C60)alkyl, (C1-C60)alkoxy, (C3-C60)cycloalkyl, (C6-C60)aryl, (C4-C60)heteroaryl, halogen, cyano, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl and tri(C6-C60)arylsilyl;

45 **[0048]** R<sub>313</sub> through R<sub>316</sub> independently represent hydrogen, (C1-C60)alkyl, (C1-C60)alkoxy, halogen, (C4-C60)heteroaryl, (C5-C60)cycloalkyl or (C6-C60)aryl, and the heteroaryl, cycloalkyl or aryl of R<sub>313</sub> through R<sub>316</sub> may be further substituted by one or more substituent(s) selected from a group consisting of (C1-C60)alkyl with or without halogen substituent(s), (C1-C60)alkoxy, (C3-C60)cycloalkyl, halogen, cyano, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl and tri(C6-C60)arylsilyl;

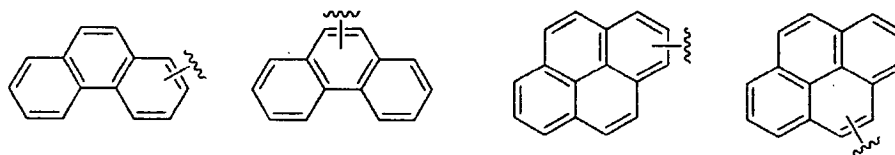
G<sub>1</sub> and G<sub>2</sub> independently represent a chemical bond or (C6-C60)arylene with or without one or more substituent(s) selected from (C1-C60)alkyl, (C1-C60)alkoxy, (C6-C60)aryl, (C4-C60)heteroaryl and halogen;

55 Ar<sub>41</sub> and Ar<sub>42</sub> represent an aryl selected from the following structures, or (C4-C60)heteroaryl:

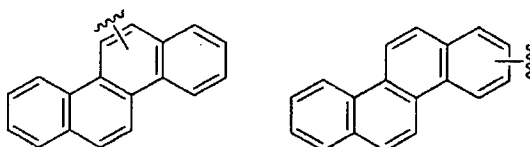
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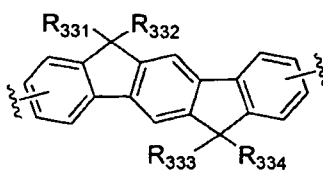
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the aryl or heteroaryl of Ar<sub>41</sub> and Ar<sub>42</sub> may be substituted by one or more substituent(s) selected from (C1-C60) alkyl, (C1-C60)alkoxy, (C6-C60)aryl and (C4-C60)heteroaryl;

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L<sub>31</sub> represents (C6-C60)arylene, (C4-C60)heteroarylene or a compound represented by the following structure:

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35

the arylene or heteroarylene of L<sub>31</sub> may be substituted by one or more substituent(s) selected from (C1-C60)alkyl, (C1-C60)alkoxy, (C6-C60)aryl, (C4-C60)heteroaryl and halogen;

R<sub>321</sub>, R<sub>322</sub>, R<sub>323</sub> and R<sub>324</sub> independently represent hydrogen, (C1-C60)alkyl or (C6-C60)aryl, or each of them may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

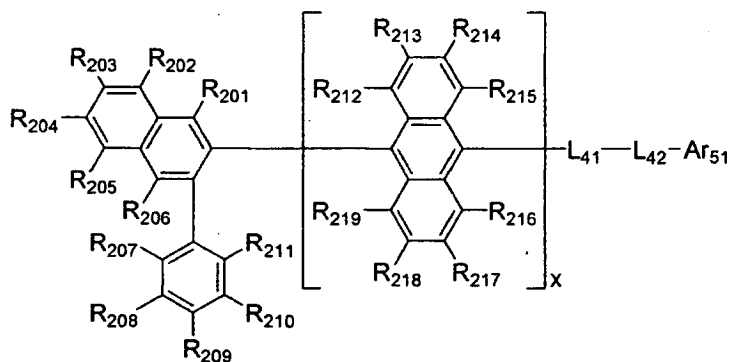
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R<sub>331</sub>, R<sub>332</sub>, R<sub>333</sub> and R<sub>334</sub> independently represent hydrogen, (C1-C60)alkyl, (C1-C60)alkoxy, (C6-C60)aryl, (C4-C60)heteroaryl or halogen, or each of them may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring.

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Chemical Formula 21

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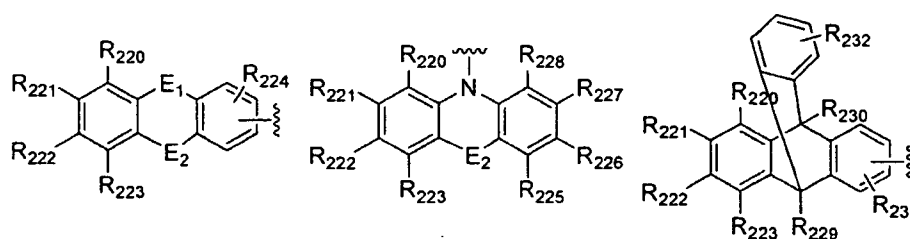
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[0049] In Chemical Formula 21,

$L_{41}$  and  $L_{42}$  independently represent a chemical bond, (C6-C60)arylene or (C3-C60)heteroarylene, and the arylene or heteroarylene of  $L_{41}$  and  $L_{42}$  may be further substituted by one or more substituent(s) selected from (C1-C60)alkyl, halogen, cyano, (C1-C60)alkoxy, (C3-C60)cycloalkyl, (C6-C60)aryl, (C3-C60)heteroaryl, tri(C1-C30)alkylsilyl, di(C1-C30)alkyl(C6-C30)arylsilyl and tri(C6-C30)arylsilyl;

$R_{201}$  through  $R_{219}$  independently represent hydrogen, halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, (C1-C60)alkoxy, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, carboxyl, nitro or hydroxyl, or each of  $R_{201}$  through  $R_{219}$  may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

$Ar_{51}$  represents (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, adamantyl, (C7-C60)bicycloalkyl, or a substituent selected from the following structures:



wherein,  $R_{220}$  through  $R_{232}$  independently represent hydrogen, halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, (C1-C60)alkoxy, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, carboxyl, nitro or hydroxyl;

$E_1$  and  $E_2$  independently represent a chemical bond,  $-(CR_{233}R_{234})_z-$ ,  $-N(R_{235})-$ ,  $-S-$ ,  $-O-$ ,  $-Si(R_{236})(R_{237})-$ ,  $-P(R_{238})-$ ,  $-C(=O)-$ ,  $-B(R_{239})-$ ,  $-In(R_{240})-$ ,  $-Se-$ ,  $-Ge(R_{241})(R_{242})-$ ,  $-Sn(R_{243})(R_{244})-$ ,  $-Ga(R_{245})-$  or  $-C(R_{246})=C(R_{247})-$ ;

$R_{233}$  through  $R_{247}$  independently represent hydrogen, halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, (C1-C60)alkoxy, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, carboxyl, nitro or hydroxyl, or each of  $R_{233}$  through  $R_{247}$  may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

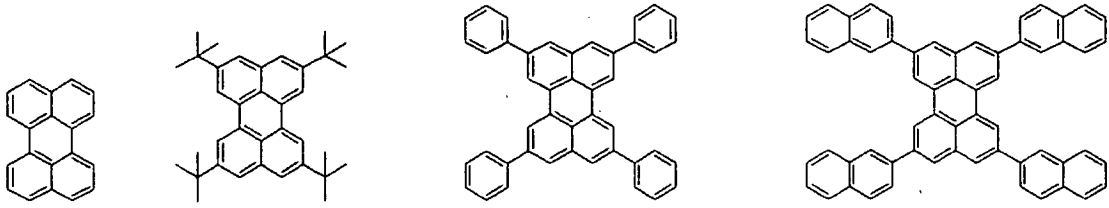
the aryl, heteroaryl, heterocycloalkyl, adamantyl or bicycloalkyl of  $Ar_{51}$ , or the alkyl, alkenyl, alkynyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, arylsilyl, alkylsilyl, alkylamino or arylamino of  $R_{201}$  through  $R_{232}$  may be further substituted by one or more substituent(s) selected from halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, (C1-C60)alkoxy, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, carboxyl, nitro and hydroxyl;

$x$  is an integer from 1 to 4; and

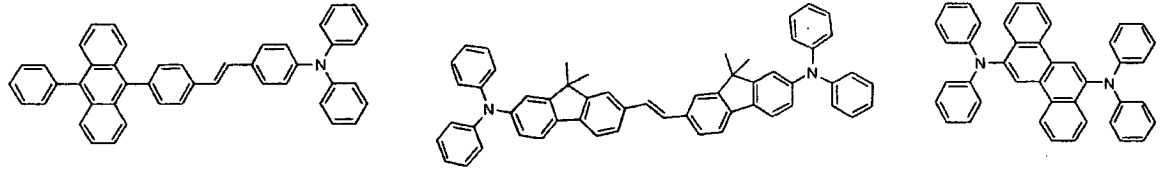
$z$  is an integer from 0 to 4.

[0050] The organic compounds and organometallic compounds with green or blue electroluminescence can be more specifically exemplified by the following compounds, but they are not restricted thereto.

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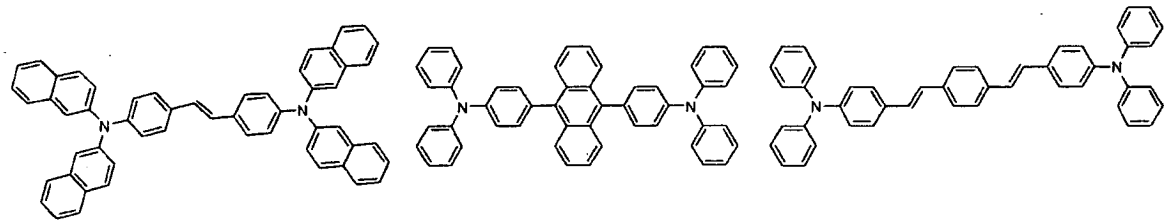


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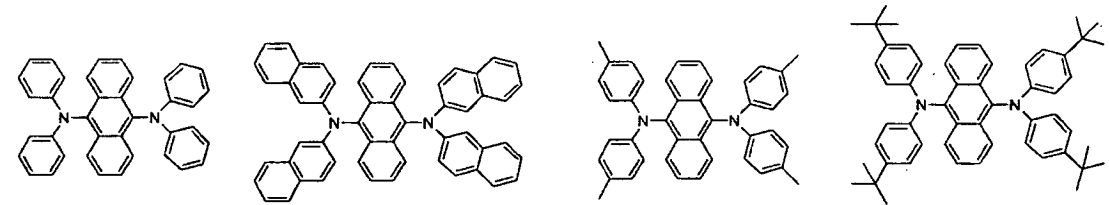
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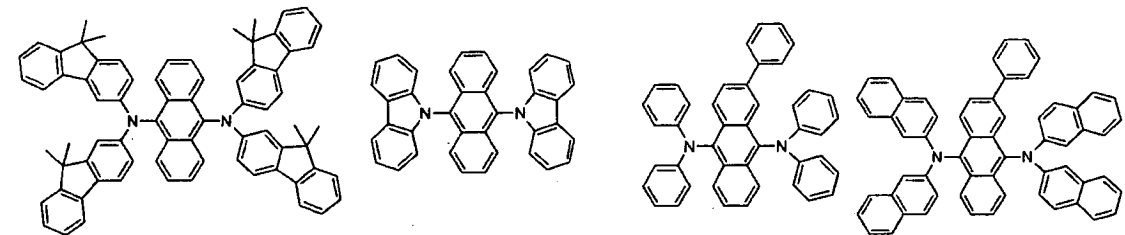
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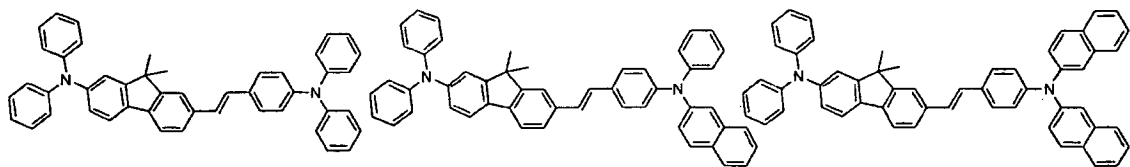
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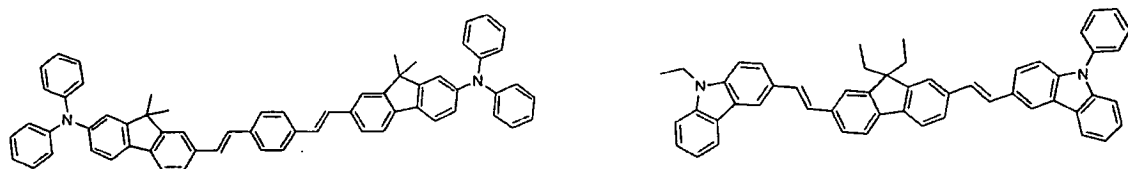
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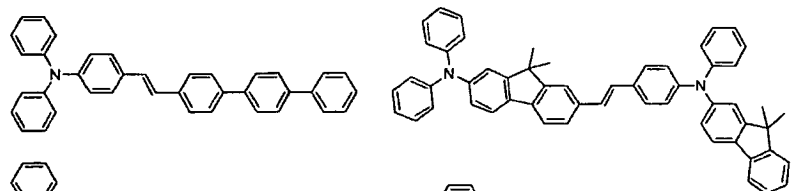


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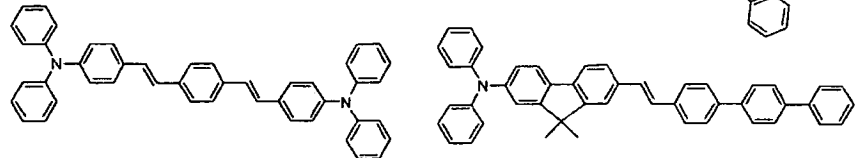
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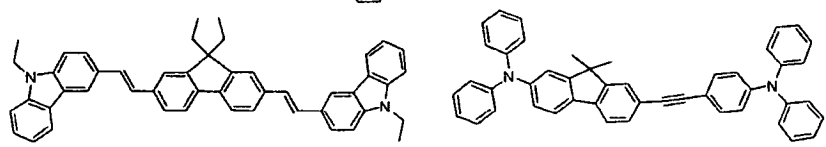
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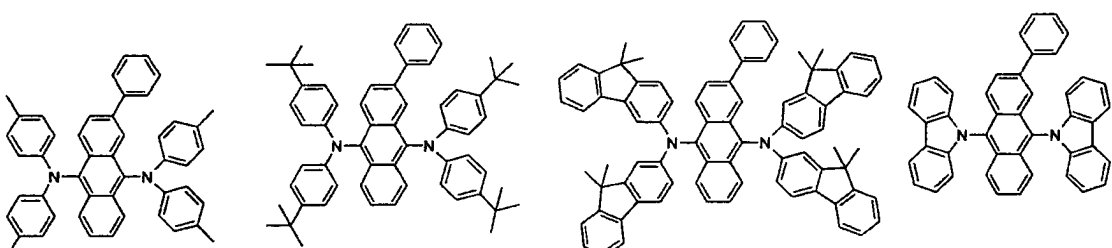


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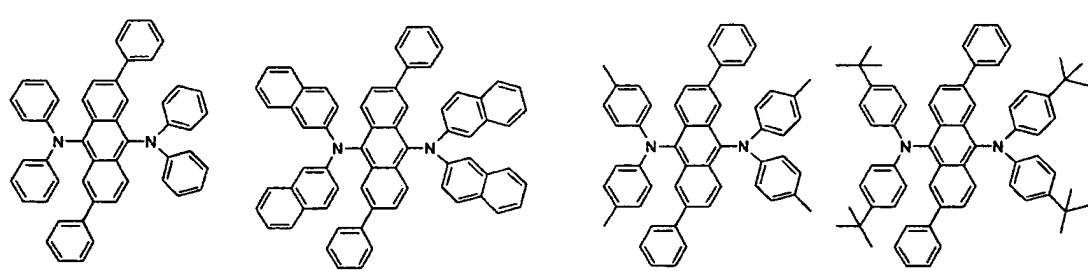
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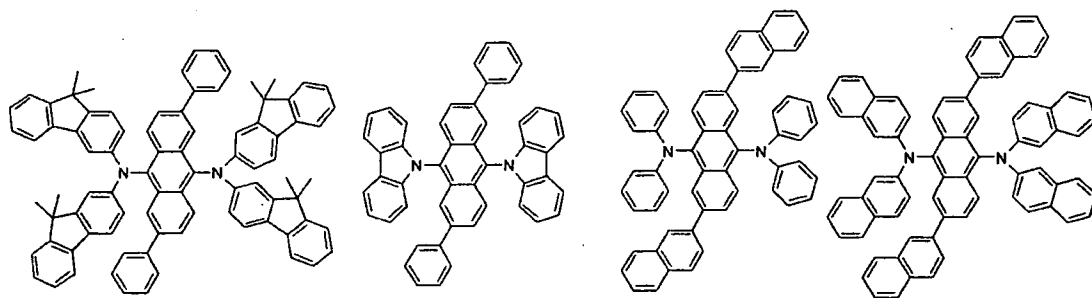


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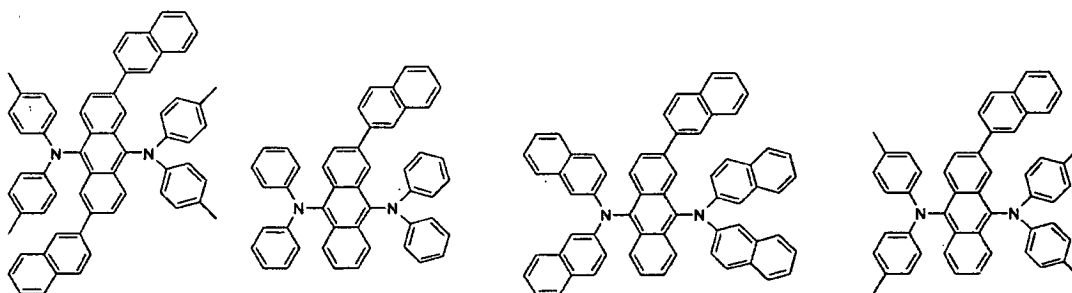
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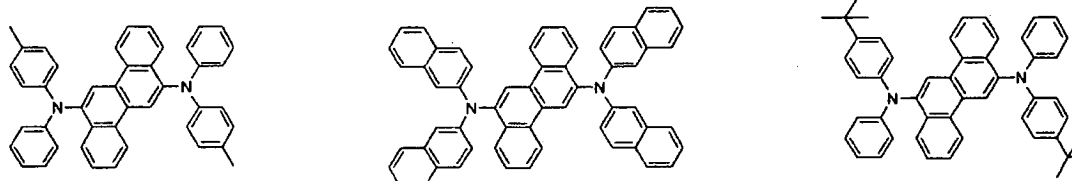
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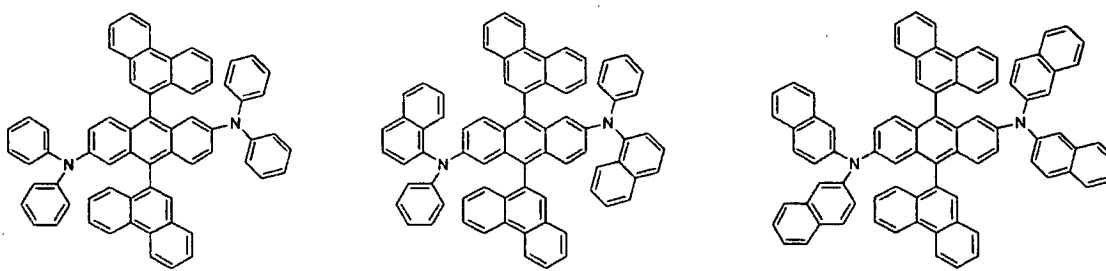
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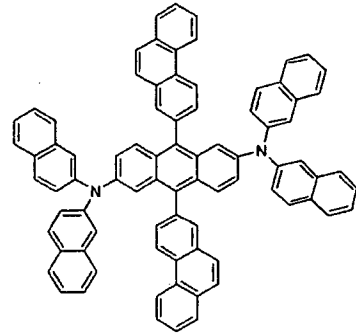
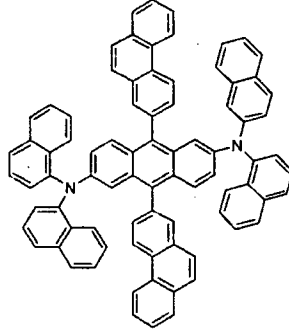
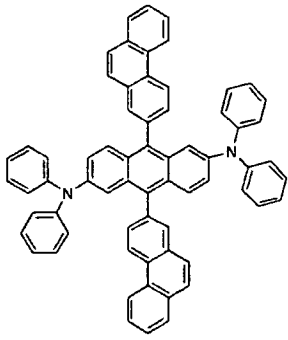


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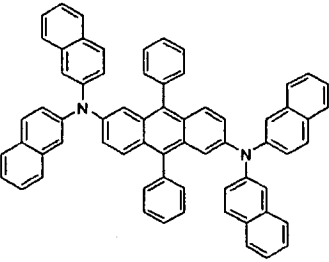
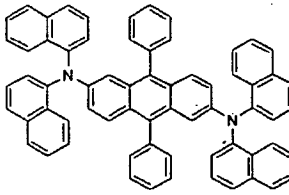
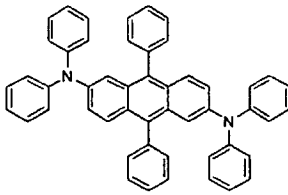
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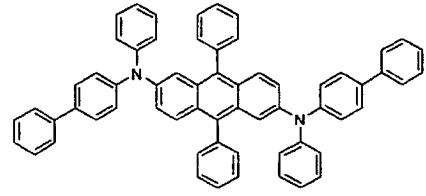
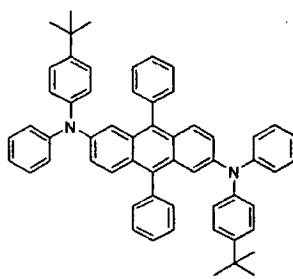
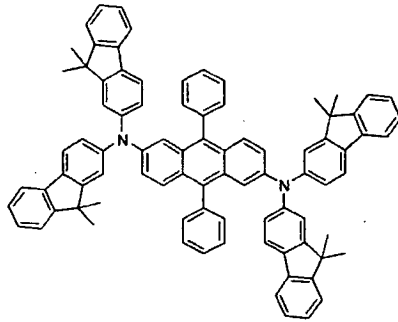
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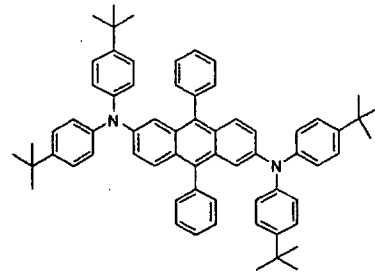
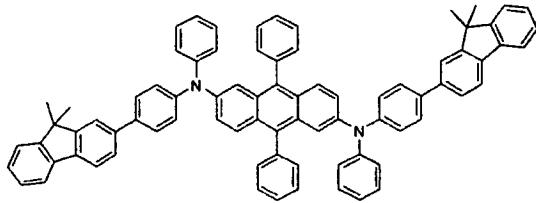
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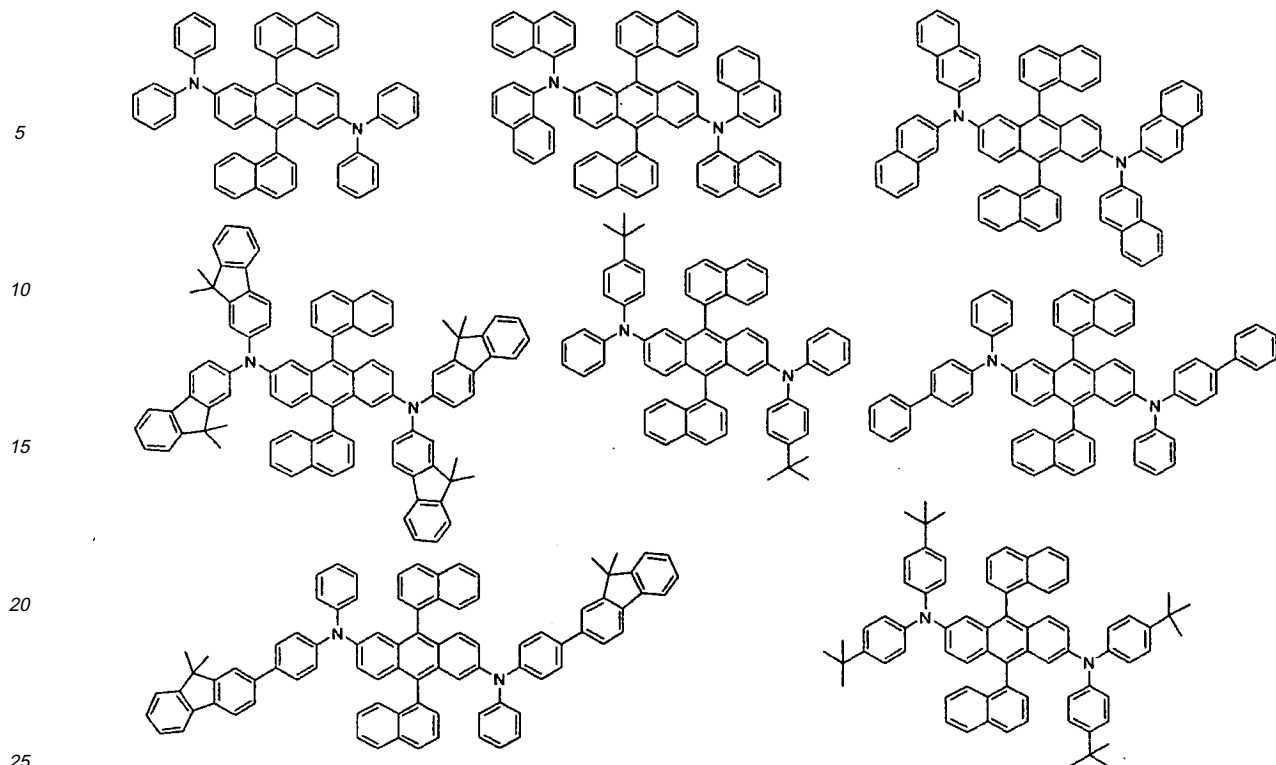


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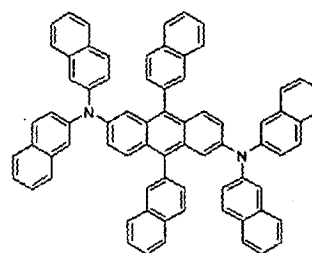
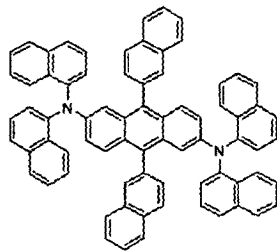
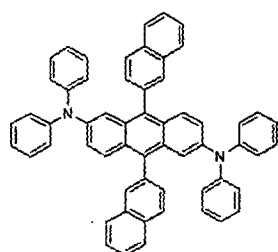
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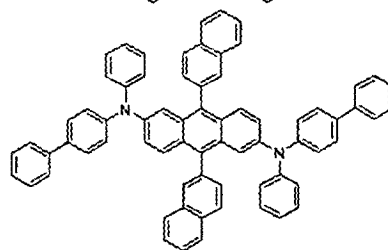
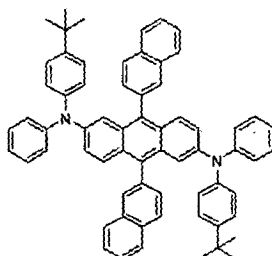
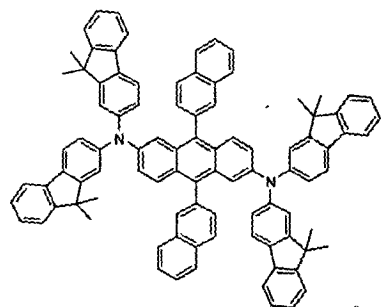
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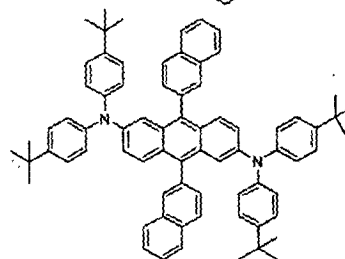
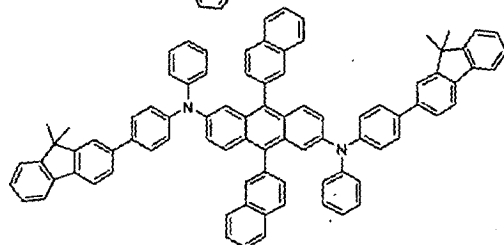
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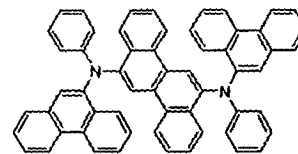
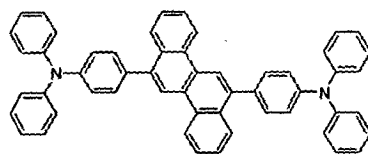
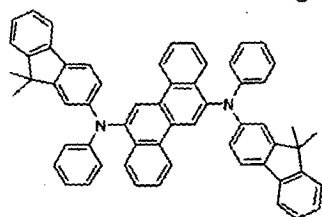
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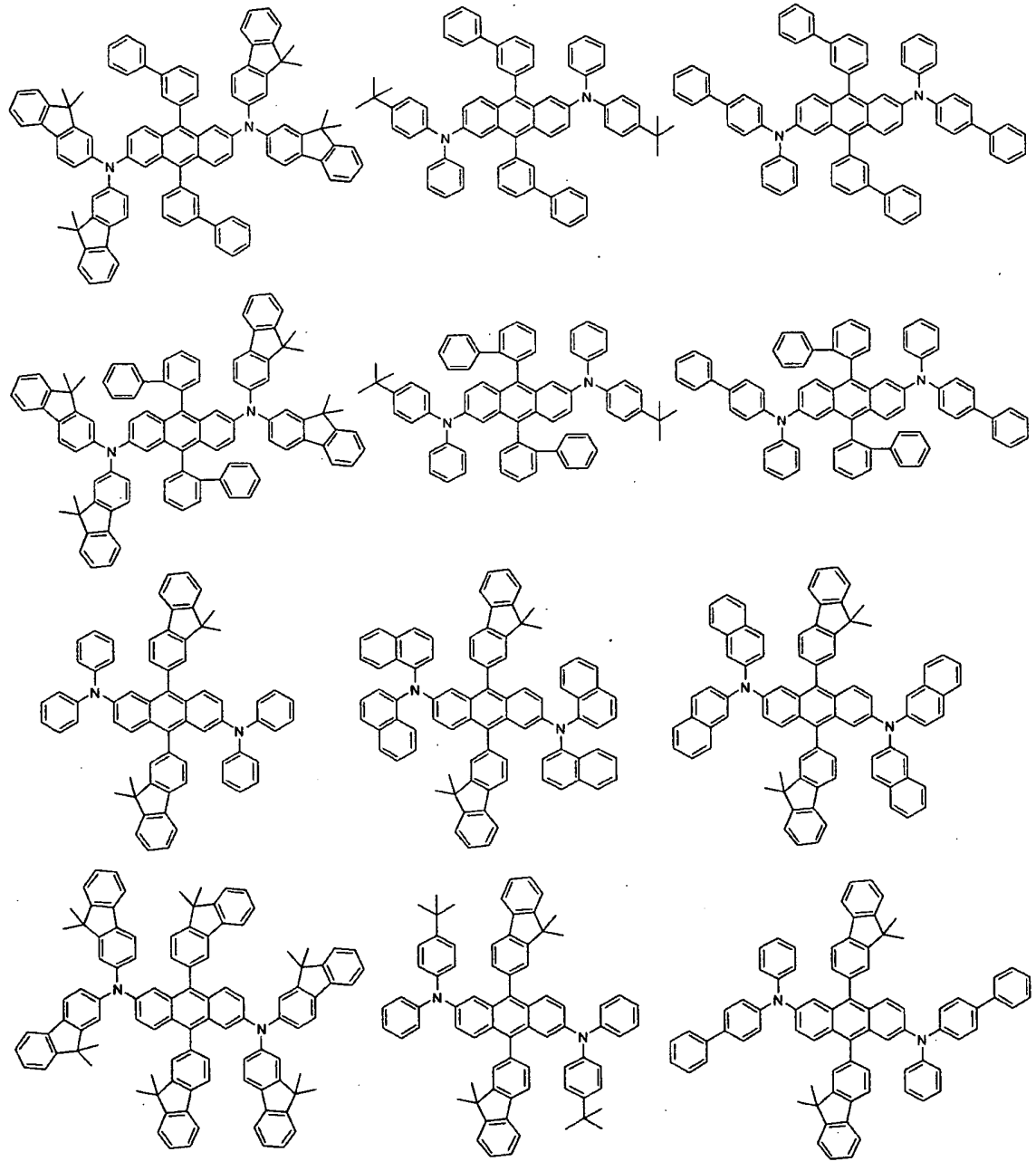
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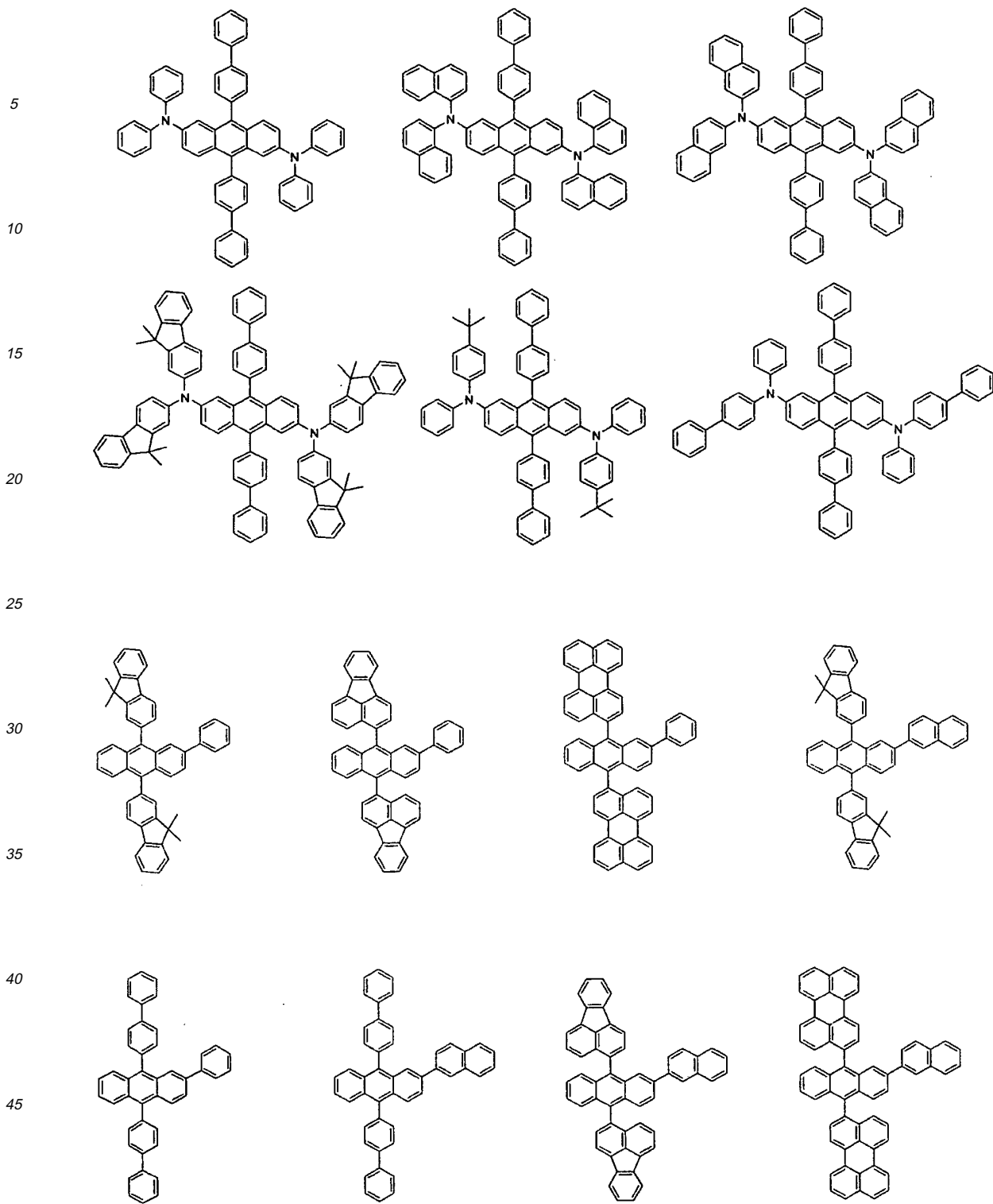
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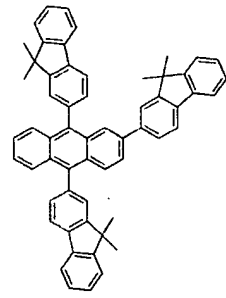
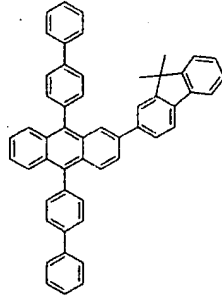
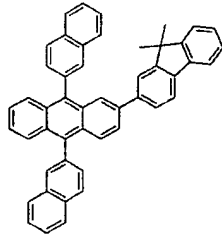
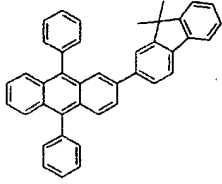
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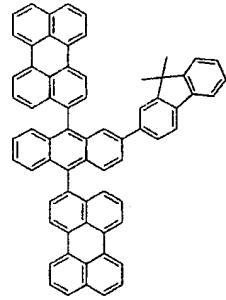
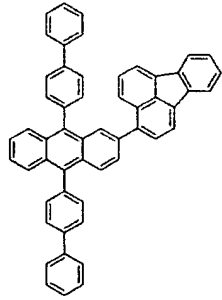
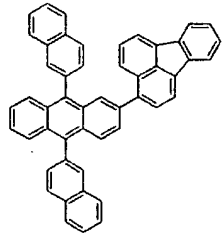
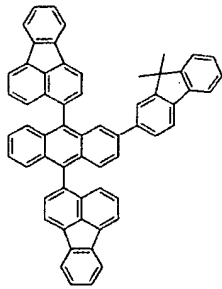


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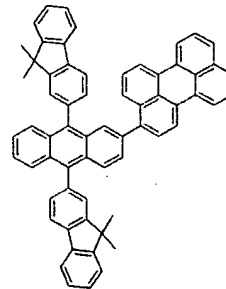
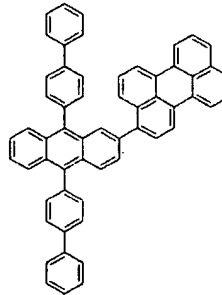
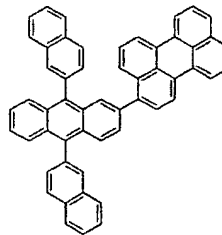
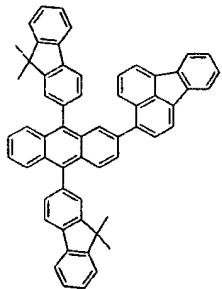
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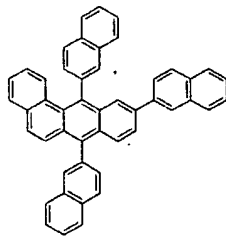
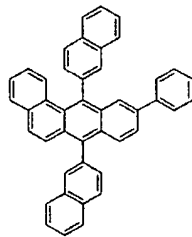
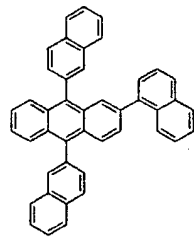
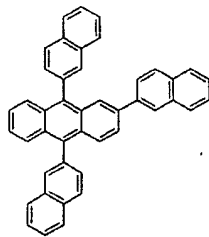
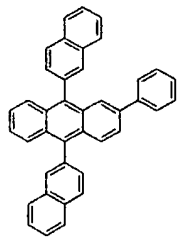
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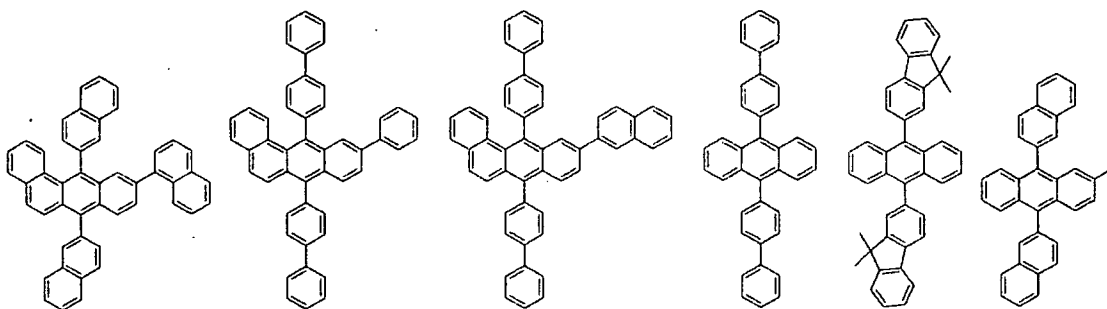


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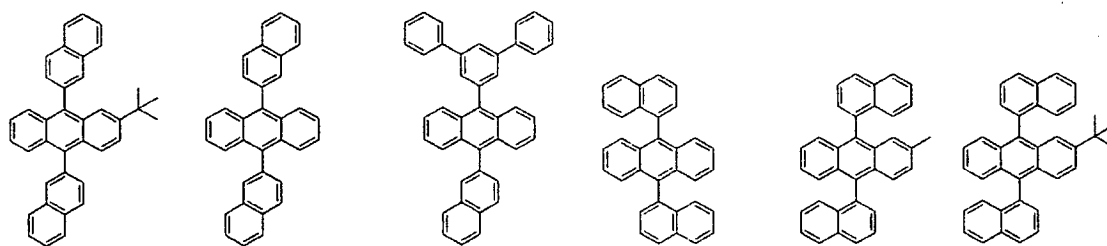
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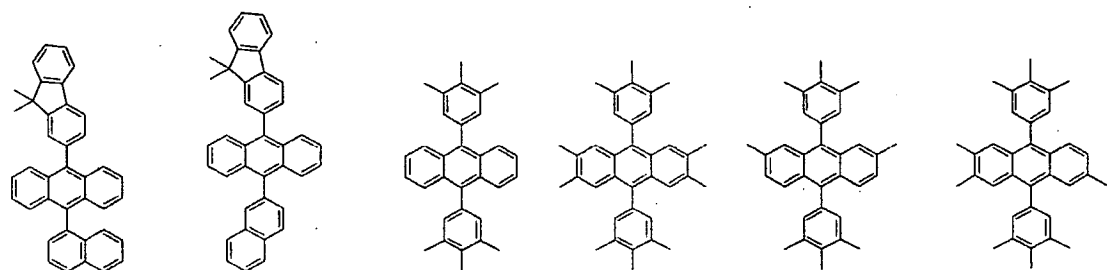
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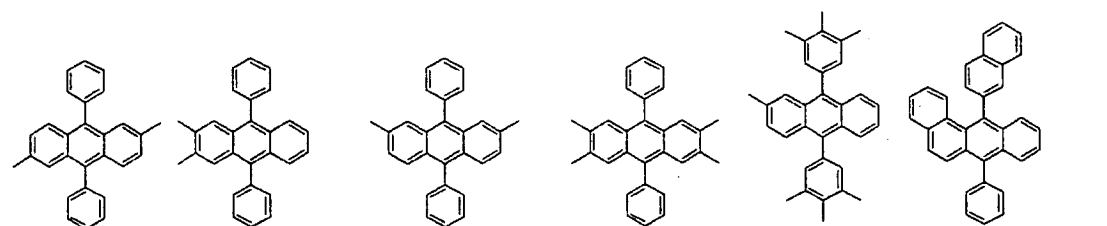
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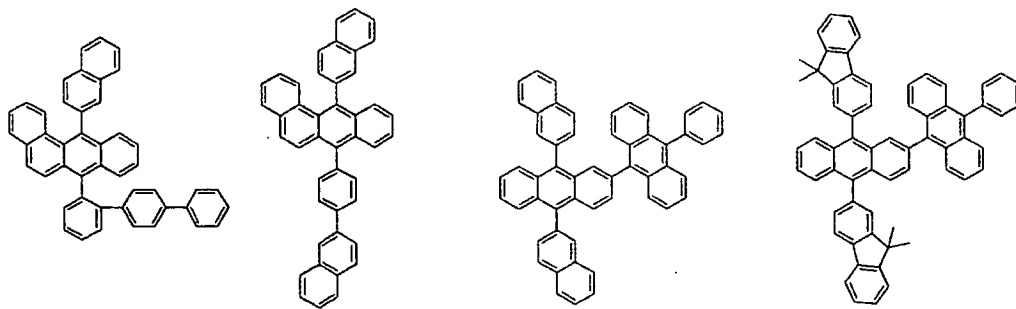


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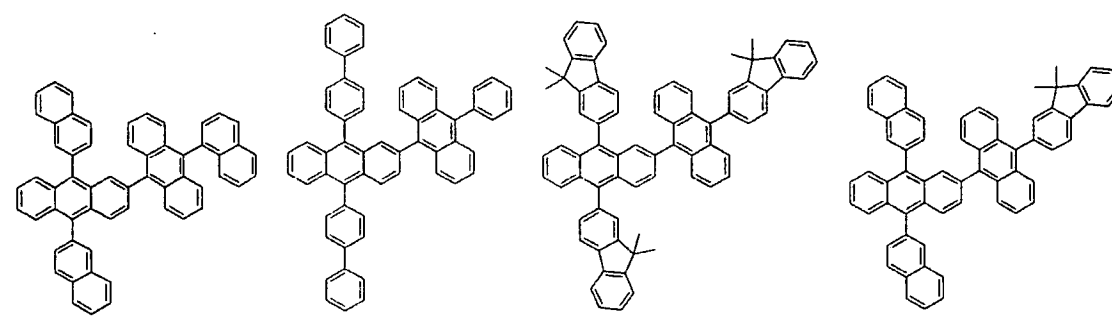
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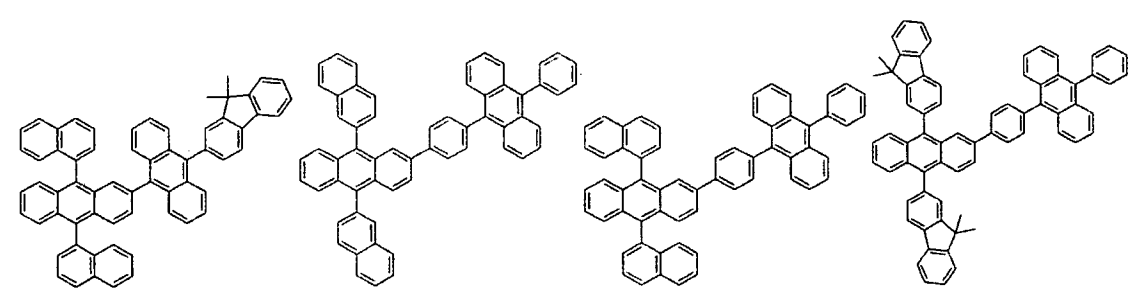
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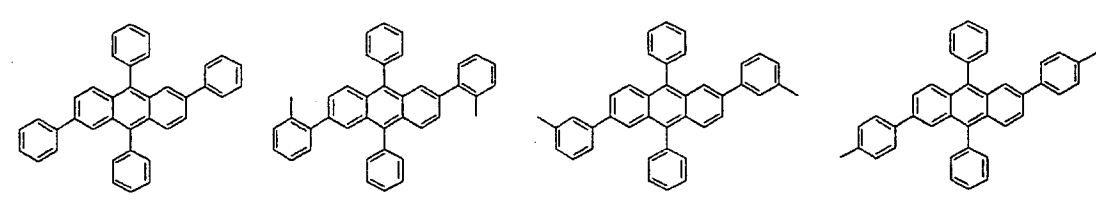
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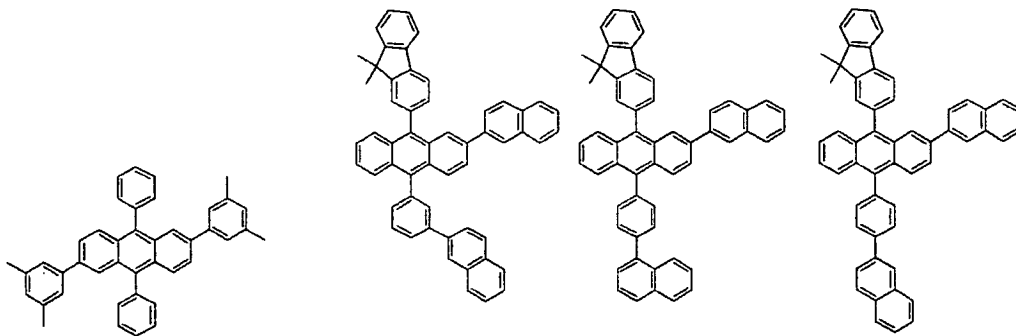


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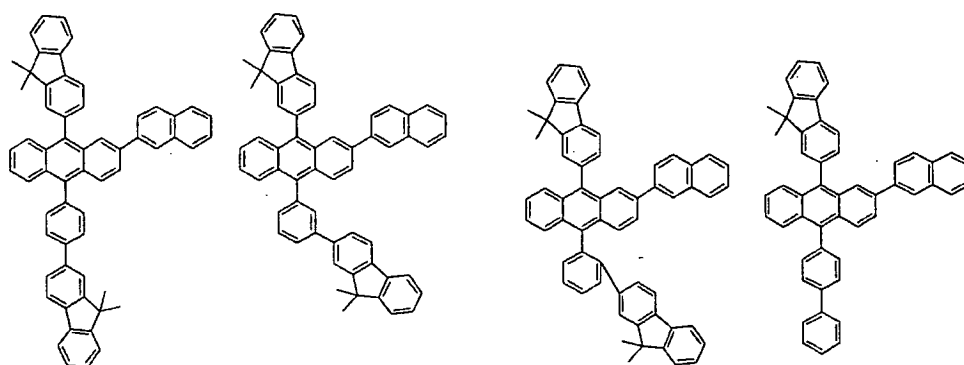
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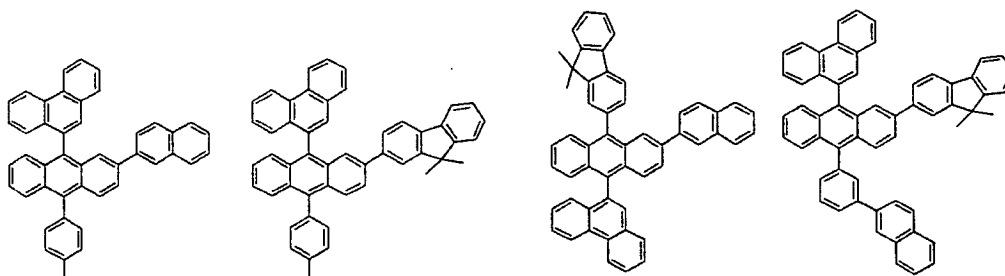
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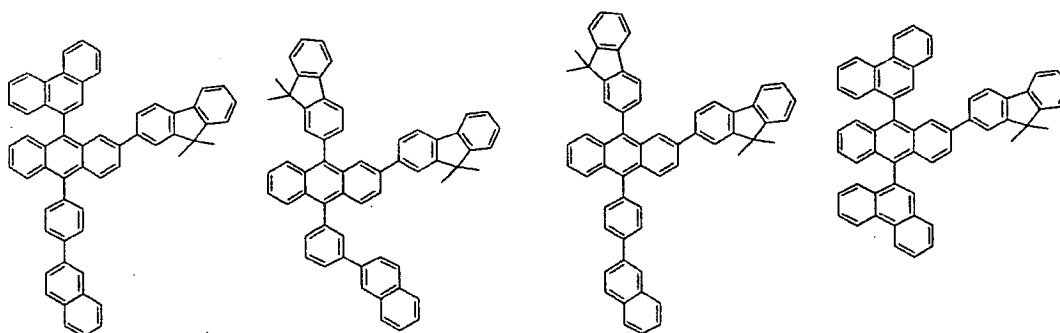
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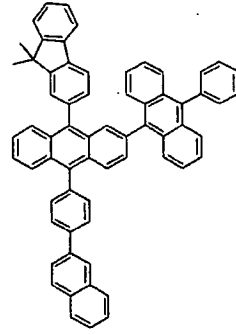
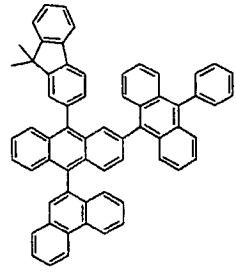
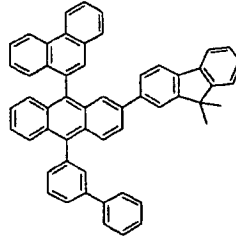
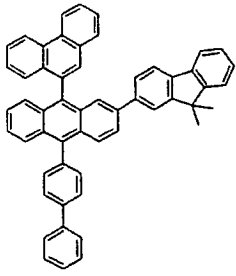
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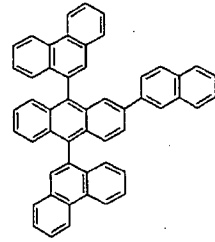
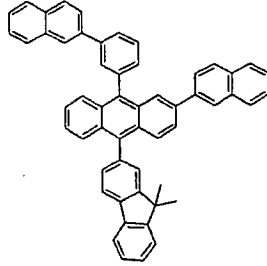
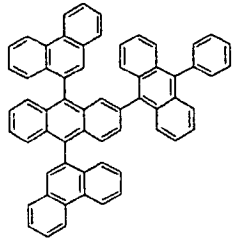
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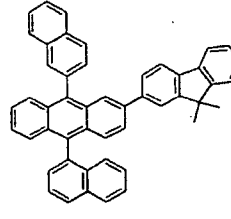
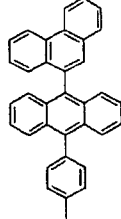
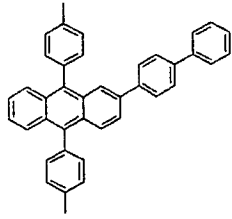
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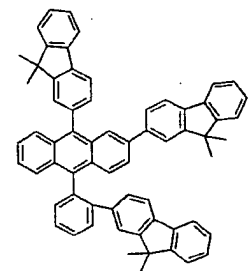
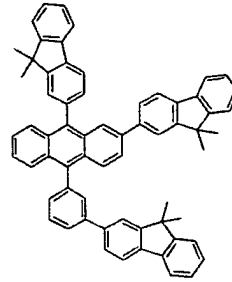
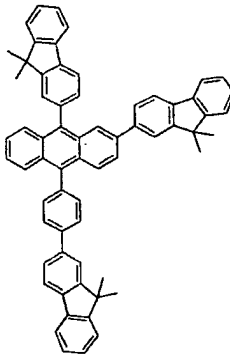
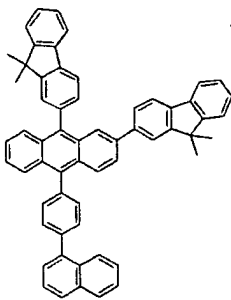
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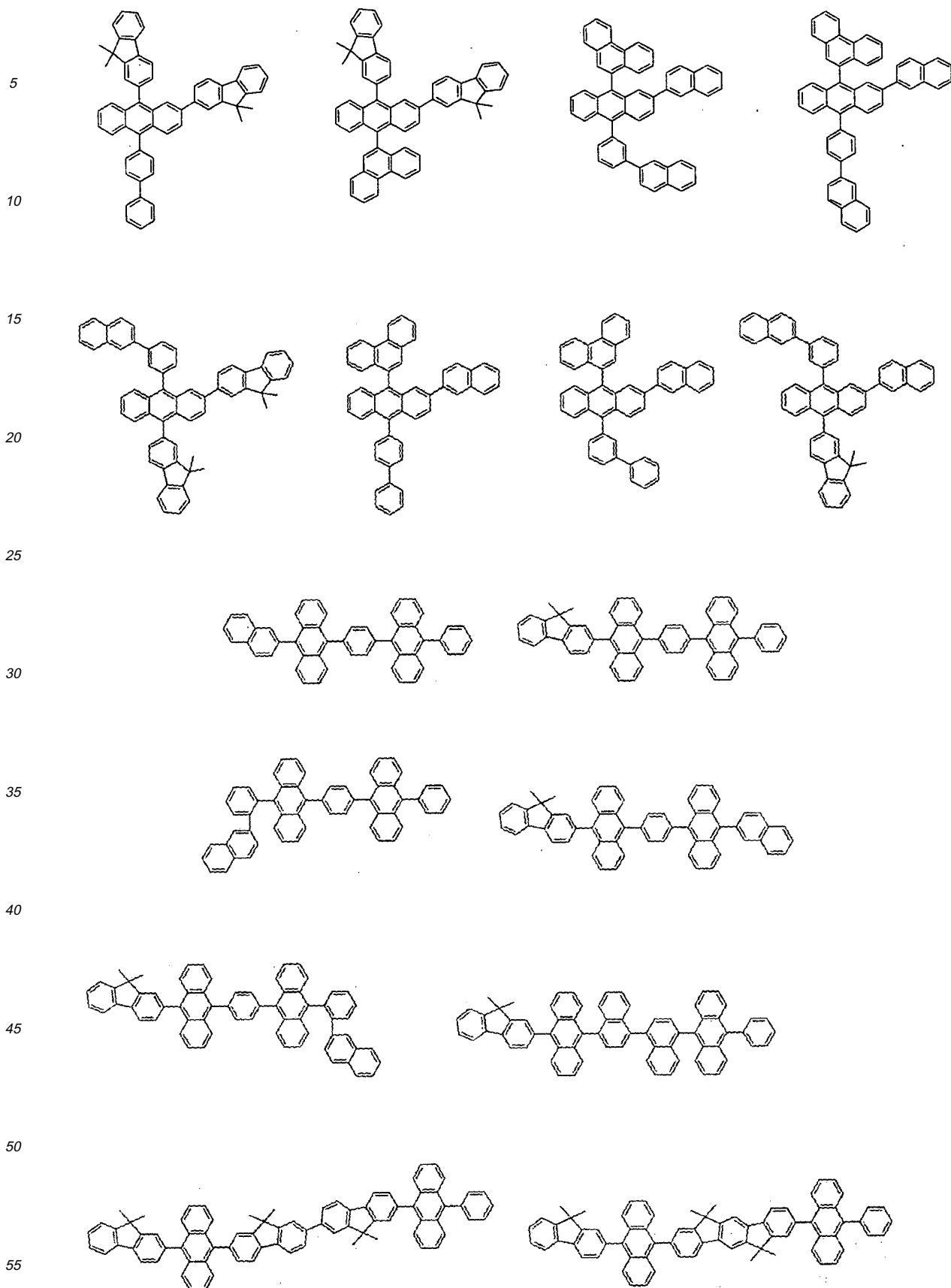
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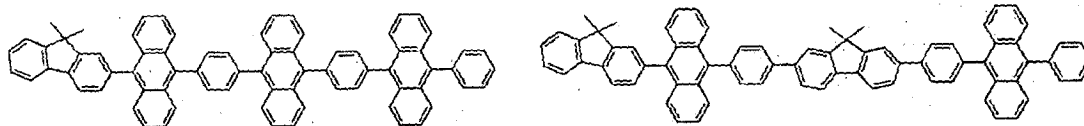
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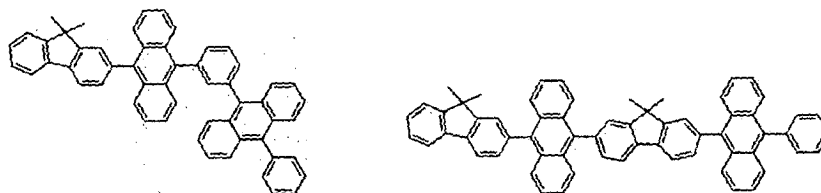
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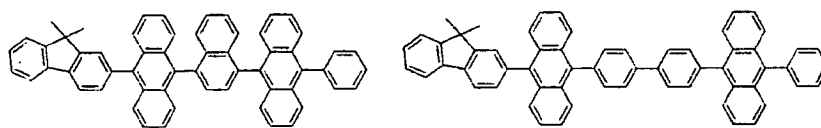


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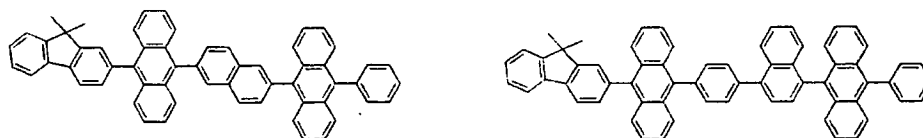


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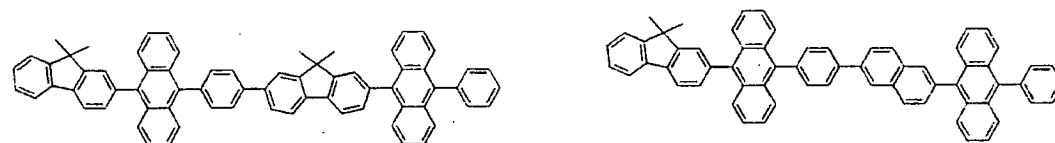


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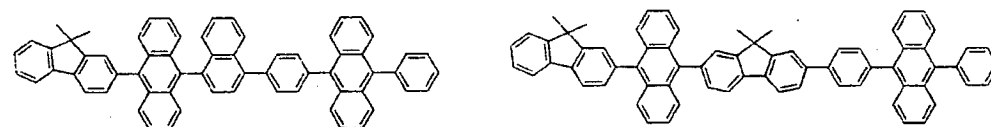
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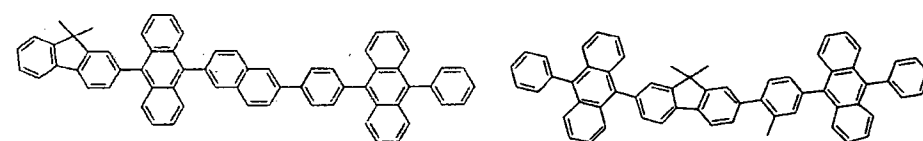


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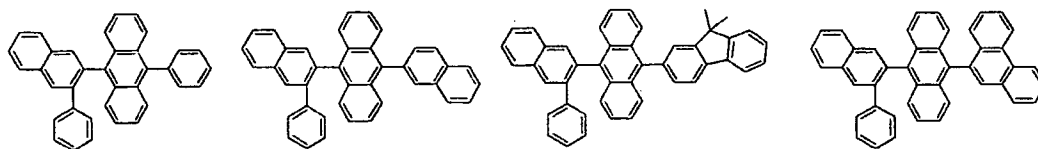


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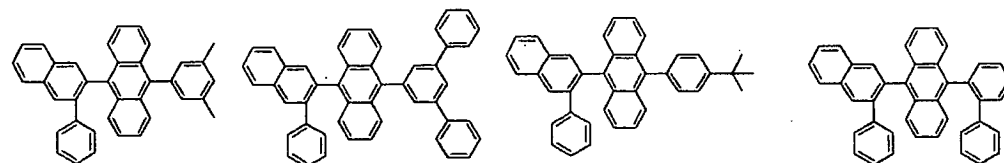


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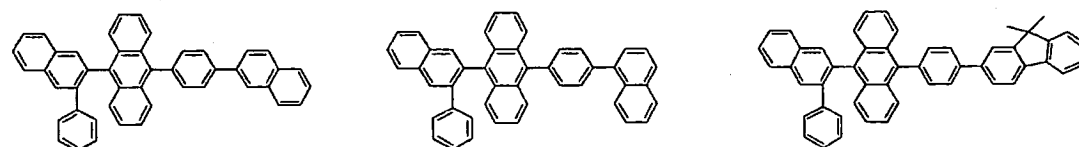


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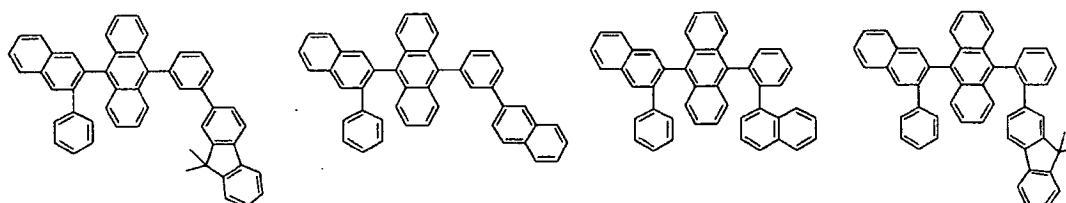
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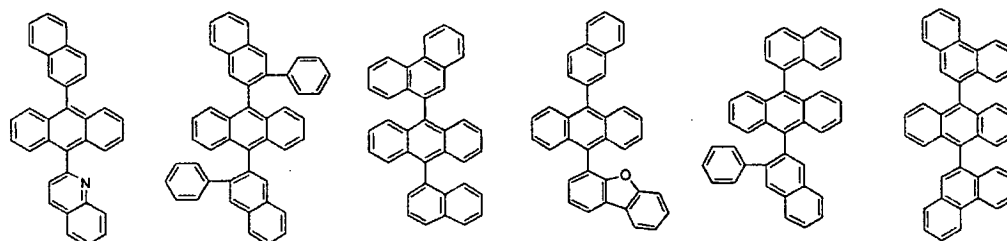
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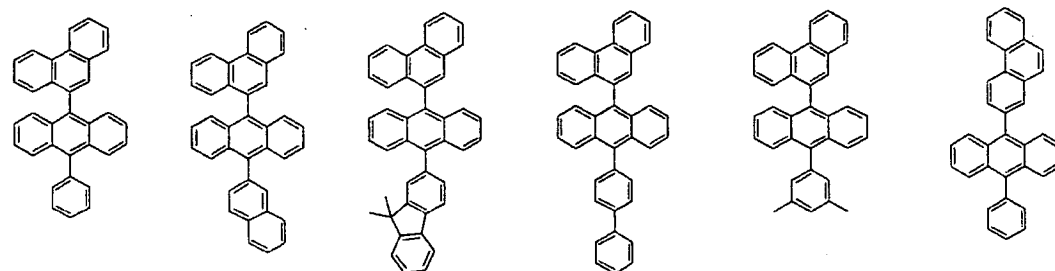
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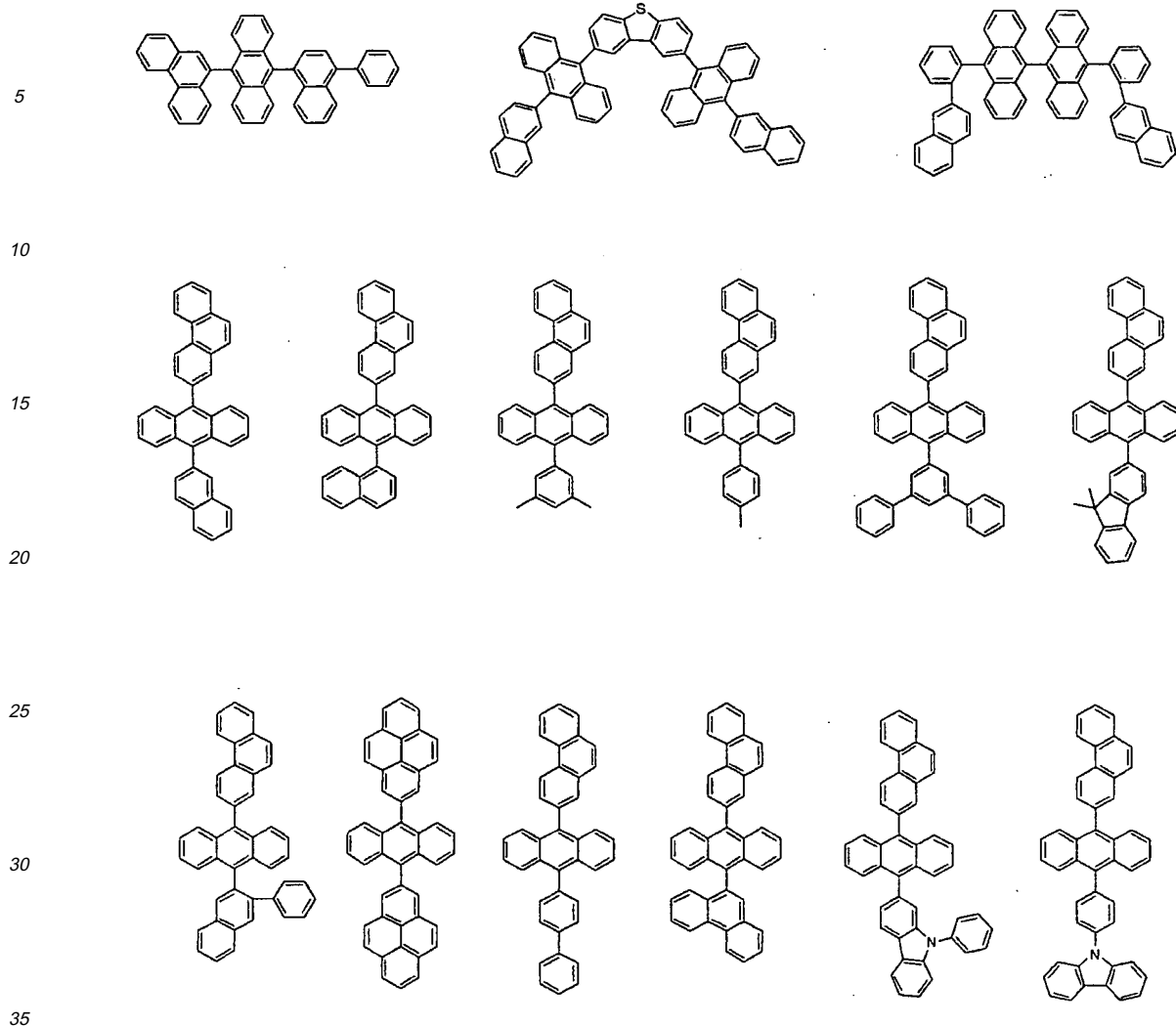


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**[0051]** In an organic electroluminescent device according to the present invention, it is preferable to displace one or more layer(s) (here-in-below, referred to as the "surface layer") selected from chalcogenide layers, metal halide layers and metal oxide layers, on the inner surface of at least one side of the pair of electrodes. Specifically, it is preferable to arrange a chalcogenide layer of silicon and aluminum metal (including oxides) on the anode surface of the EL medium layer, and a metal halide layer or a metal oxide layer on the cathode surface of the EL medium layer. As the result, stability in operation can be obtained.

**[0052]** Examples of chalcogenides preferably include  $\text{SiO}_x$  ( $1 \leq x \leq 2$ ),  $\text{AlO}_x$  ( $1 \leq x \leq 1.5$ ),  $\text{SiON}$ ,  $\text{SiAlON}$ , or the like. Examples of metal halides preferably include  $\text{LiF}$ ,  $\text{MgF}_2$ ,  $\text{CaF}_2$ , fluorides of lanthanides or the like. Examples of metal oxides preferably include  $\text{CS}_2\text{O}$ ,  $\text{Li}_2\text{O}$ ,  $\text{MgO}$ ,  $\text{SrO}$ ,  $\text{BaO}$ ,  $\text{CaO}$ , or the like.

**[0053]** In an organic electroluminescent device according to the present invention, it is also preferable to arrange, on at least one surface of the pair of electrodes thus manufactured, a mixed region of electron transport compound and a reductive dopant, or a mixed region of a hole transport compound with an oxidative dopant. Accordingly, the electron transport compound is reduced to an anion, so that injection and transportation of electrons from the mixed region to an EL medium are facilitated. In addition, since the hole transport compound is oxidized to form a cation, injection and transportation of holes from the mixed region to an EL medium are facilitated. Preferable oxidative dopants include various Lewis acids and acceptor compounds. Preferable reductive dopants include alkali metals, alkali metal compounds, alkaline earth metals, rare-earth metals, and mixtures thereof.

**[0054]** The organic electroluminescent compounds according to the invention, having a backbone of more excellent EL properties and thermal stability than conventional phosphorescent materials, provide higher quantum efficiency and lower operation voltage as compared to conventional materials. Thus, if an organic electroluminescent compound according to the present invention is applied to an OLED panel, further enhanced results are anticipated in development

of OLED's having medium to large size. If the compound is applied to an organic solar cell as a material of high efficiency, more excellent properties are anticipated as compared to conventional materials.

Best Mode

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**[0055]** The present invention is further described with respect to the processes for preparing the novel compounds according to the present invention, which are provided for better understanding of the present invention only but are not intended to limit the scope of the invention by any means.

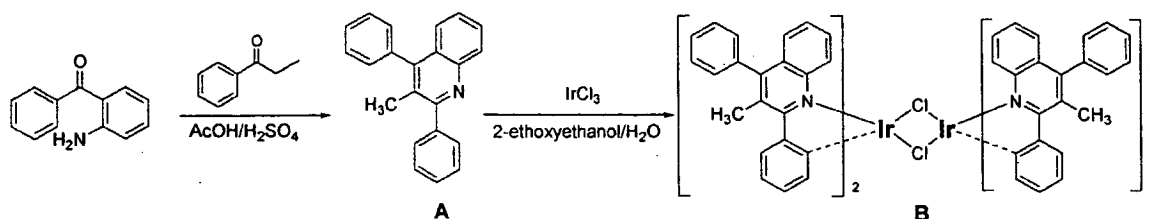
10 Preparation Examples

[Preparation Example 1] Preparation of Compound (993)

**[0056]**

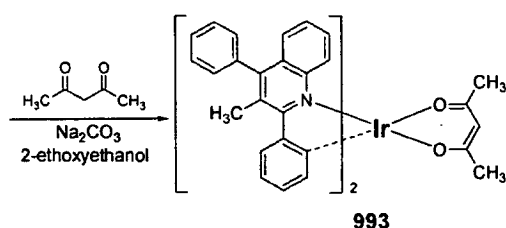
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#### Preparation of Compound (A)

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**[0057]** A reaction vessel was charged with 2-aminobenzophenone (10.0 g, 50.7 mmol), 1-methylacetophenone (6.7 mL, 50.7 mmol), glacial acetic acid (169 mL) and sulfuric acid (10 drops), and the mixture was stirred under reflux in the presence of argon gas atmosphere. When the reaction was completed, the mixture was cooled to room temperature. An excess amount of aqueous ammonium hydroxide solution was added thereto, and the solid produced was filtered and washed with distilled water. Purification via column chromatography gave Compound (A) (8.9 g, 60%).

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#### Preparation of Compound (B)

**[0058]** A reaction vessel was charged with Compound (A) (7.0 g, 23.7 mmol), iridium chloride (3.2 g, 10.8 mmol), 2-ethoxyethanol (36 mL) and distilled water (12 mL), and the mixture was heated under reflux in the presence of argon gas atmosphere for 24 hours. When the reaction was completed, the reaction mixture was cooled to room temperature. The precipitate was filtered and thoroughly dried to obtain Compound (B) (4.0 g, 71%).

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#### Preparation of Compound (993)

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**[0059]** Compound (B) (4.0 g, 3.9 mmol), 2,4-pentanedione (0.79 mL, 7.7 mmol) and sodium carbonate (2.1 g, 19.3 mmol) were dissolved in 2-ethoxyethanol (100 mL), and the solution was heated for 4 hours. When the reaction was completed, the reaction mixture was cooled to room temperature, and the solid precipitate produced was filtered. Purification via silica gel column chromatography and recrystallization gave Compound (993) (1.7 g, 25%) as red crystal.

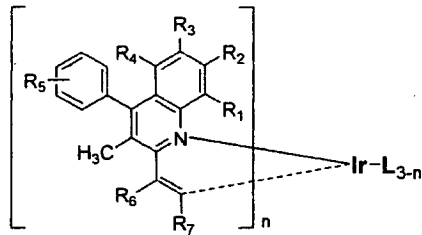
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**[0060]** According to the same procedure as Preparation Example 1, organic phosphorescent compounds (Compound 1 through Compound 1004) in Table 1 were prepared, of which the  $^1\text{H}$  NMR and MS/FAB data are listed in Table 2.

Table 1

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
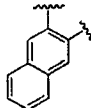
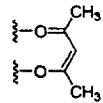


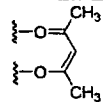

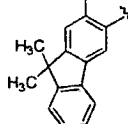
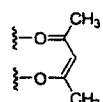

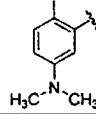
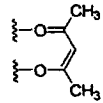

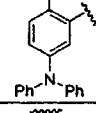
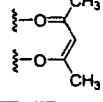


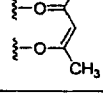

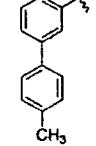
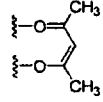

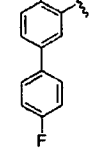
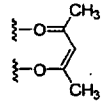

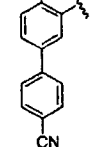
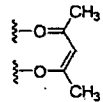
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Comp. No.	R <sub>1</sub>	R <sub>2</sub>	R <sub>3</sub>	R <sub>4</sub>	R <sub>5</sub>	R <sub>6</sub> , R <sub>7</sub>	L	n
1	H	H	H	H	Ph	CH <sub>3</sub>	CH <sub>3</sub>	2
2	H	H	H	H	Ph	F	CH <sub>3</sub>	2
3	H	H	H	H	Ph	CN	CH <sub>3</sub>	2
4	H	H	H	H	Ph	C(CH <sub>3</sub> ) <sub>3</sub>	CH <sub>3</sub>	2
5	H	H	H	H	Ph	Si(CH <sub>3</sub> ) <sub>3</sub>	CH <sub>3</sub>	2
6	H	H	H	H	Ph	SiPh <sub>3</sub>	CH <sub>3</sub>	2
7	H	H	H	H	Ph	OCH <sub>3</sub>	CH <sub>3</sub>	2
8	H	H	H	H	Ph	CF <sub>3</sub>	CH <sub>3</sub>	2
9	H	H	H	H	Ph	COCH <sub>3</sub>	CH <sub>3</sub>	2

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5	10	H	H	H	H				2
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15	12	H	H	H	H				2
20	13	H	H	H	H				2
25	14	H	H	H	H				2
30	15	H	H	H	H				2
35	16	H	H	H	H				2
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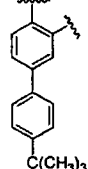
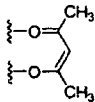

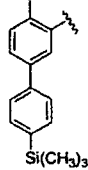
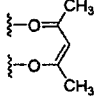

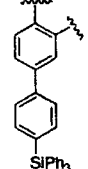
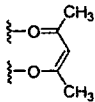

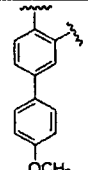
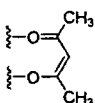

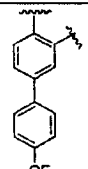
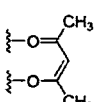

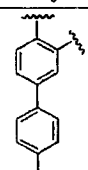
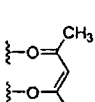

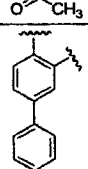
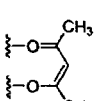
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
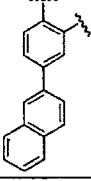
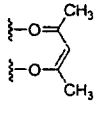

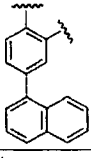
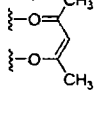

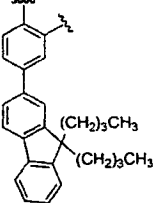
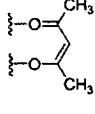

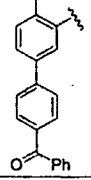
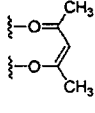

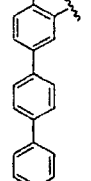
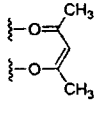

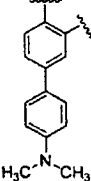
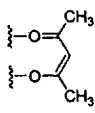

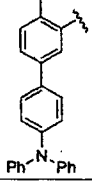
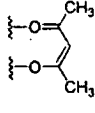
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
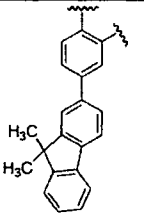
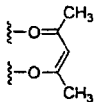

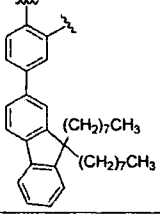
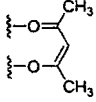

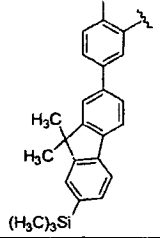
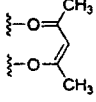

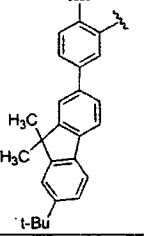
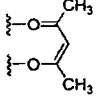

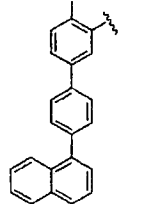
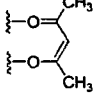
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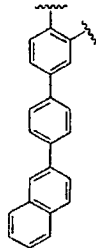
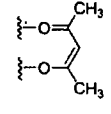

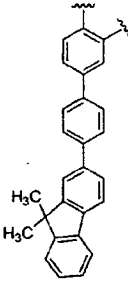
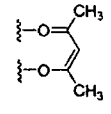

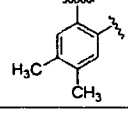
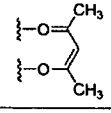

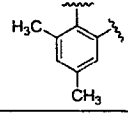
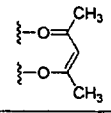

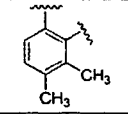
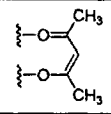

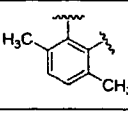
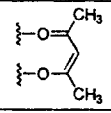

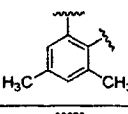
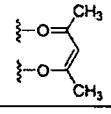

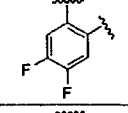
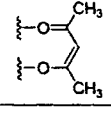

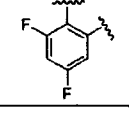
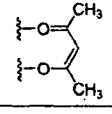
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

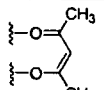


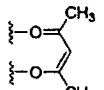


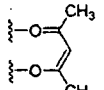


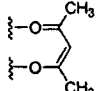


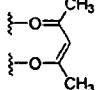


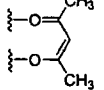


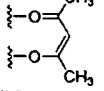


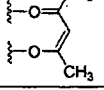


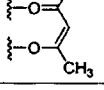


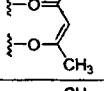


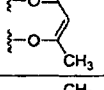


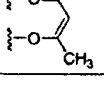
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
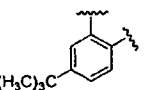


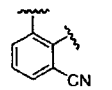


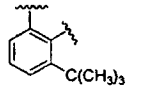


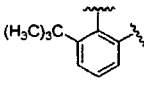


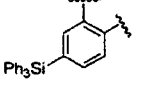


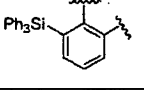


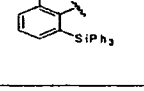


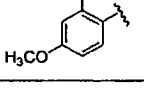


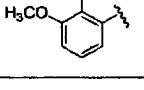


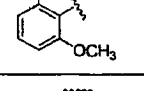


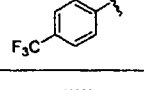


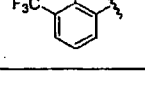

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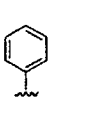
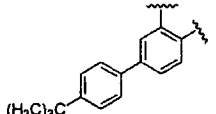
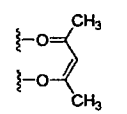
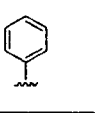
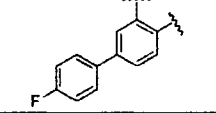
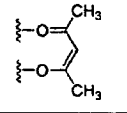
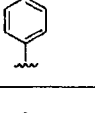
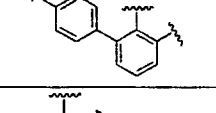
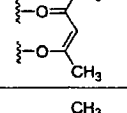
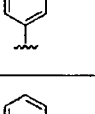
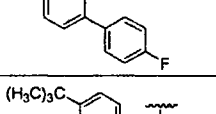
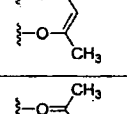
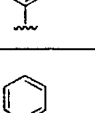
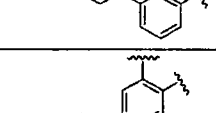
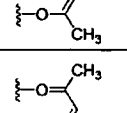
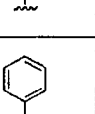
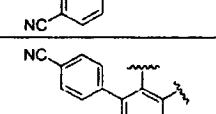
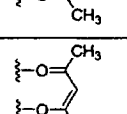
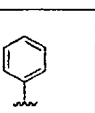
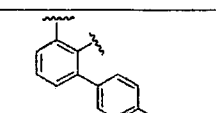
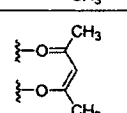
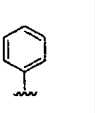
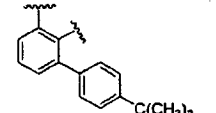
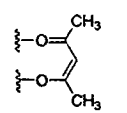
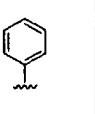
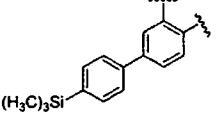
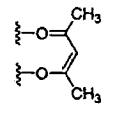
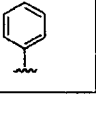
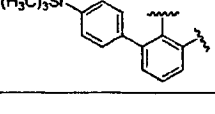
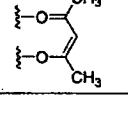



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
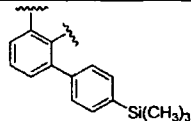
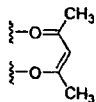

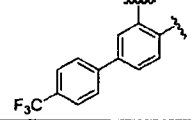
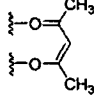

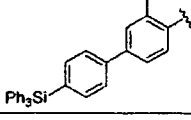
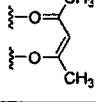

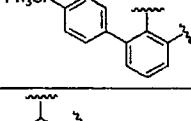
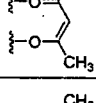

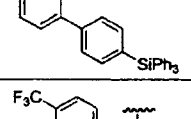
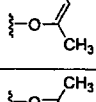

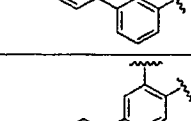
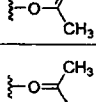

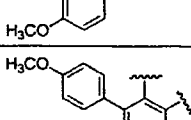
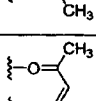

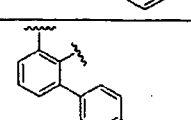
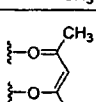

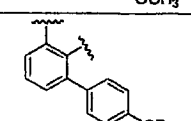
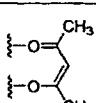

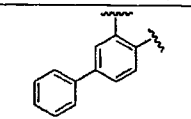
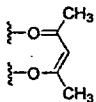



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
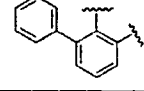


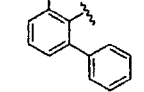


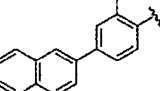


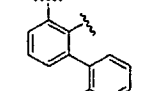


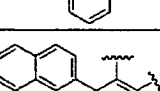


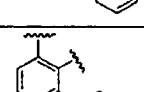


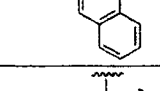


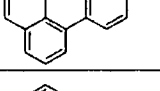


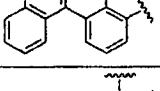


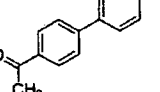

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
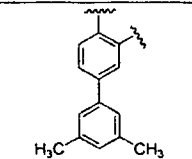
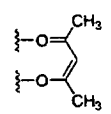

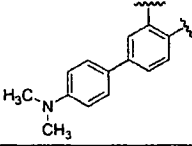
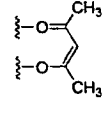

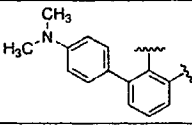
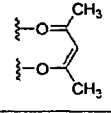

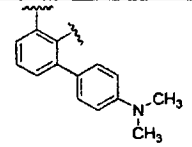
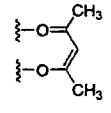

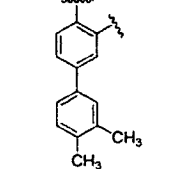
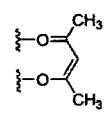

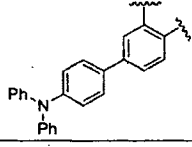
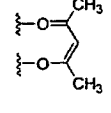

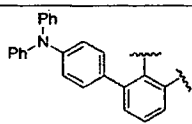
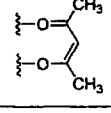

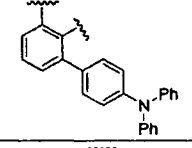
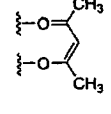

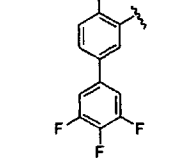
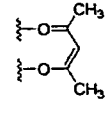
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
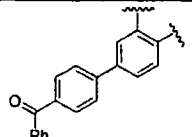
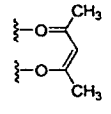

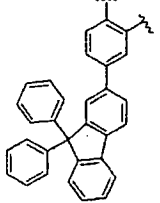
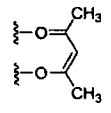

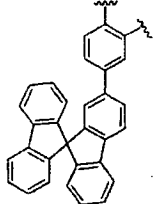
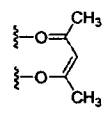

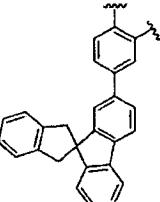
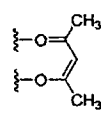

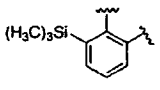
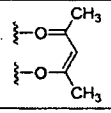

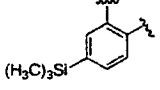
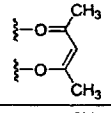

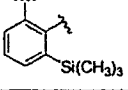
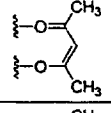

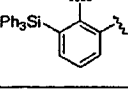
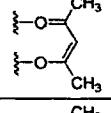

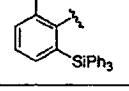
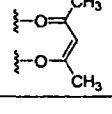
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
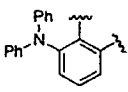
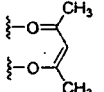

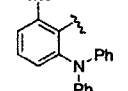
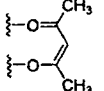

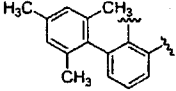
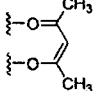

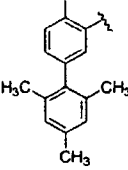
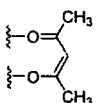

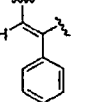
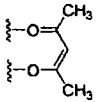

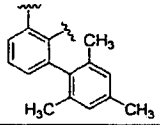
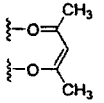

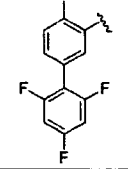
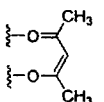

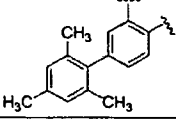
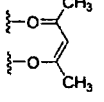

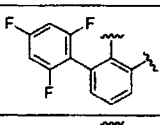
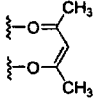

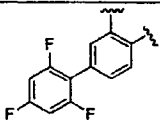
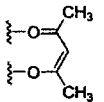
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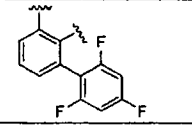
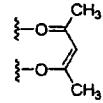

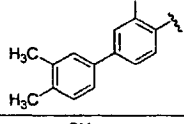
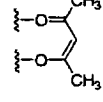

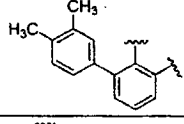
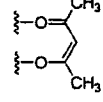

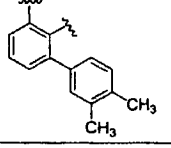
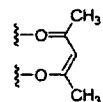

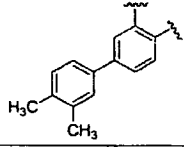
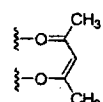

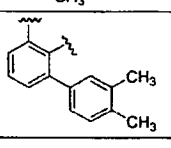
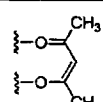

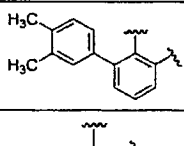
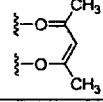

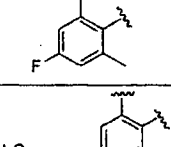
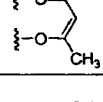

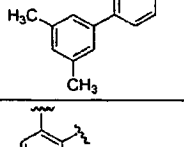
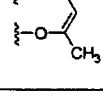

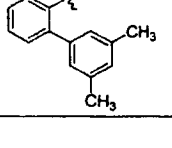
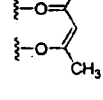
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
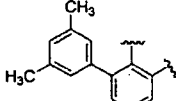
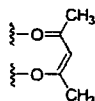

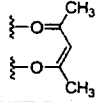

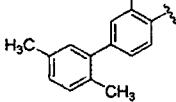
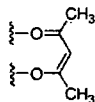

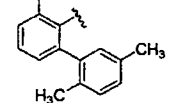
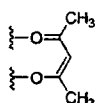

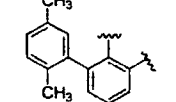
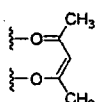

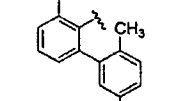
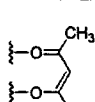

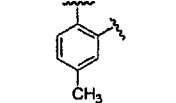
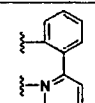

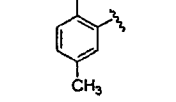
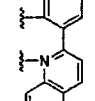

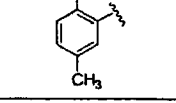
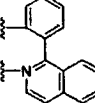

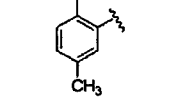
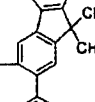
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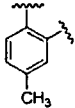
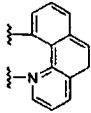

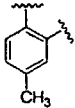
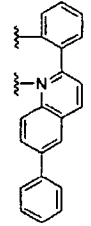

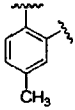
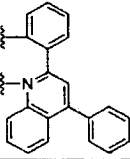

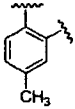
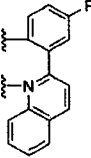

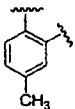
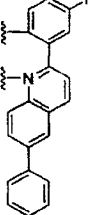

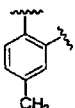
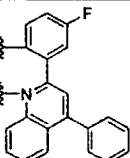

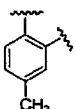
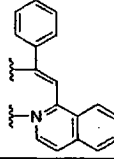

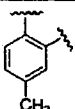
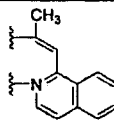
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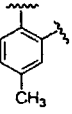
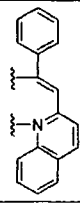

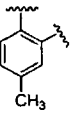
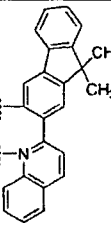

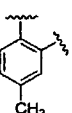
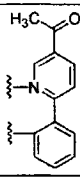

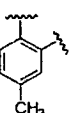
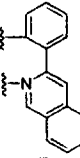

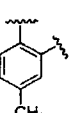
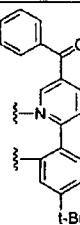

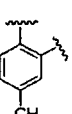
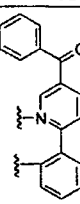

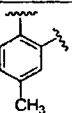
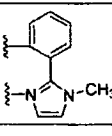
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
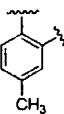
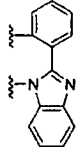

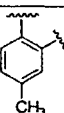


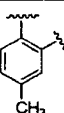
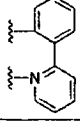

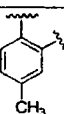

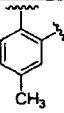


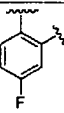


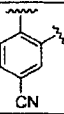


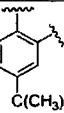


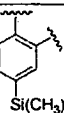


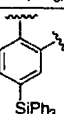


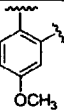

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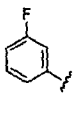
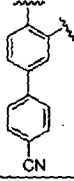
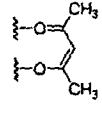
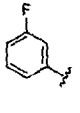
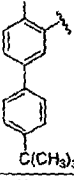
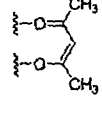
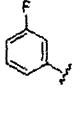
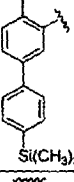
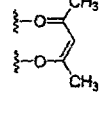
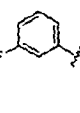
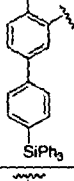
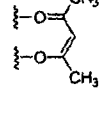
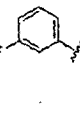
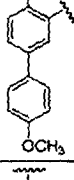
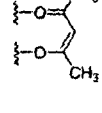
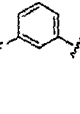
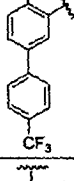
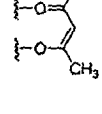
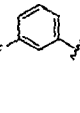
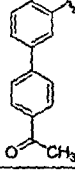
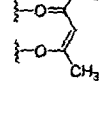
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
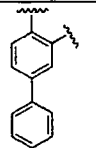
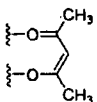

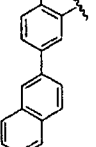
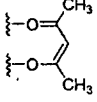

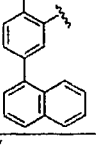
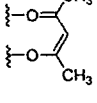

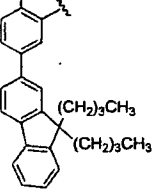
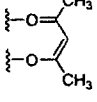

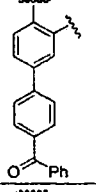
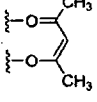

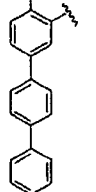
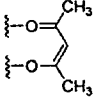

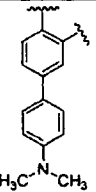
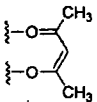
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
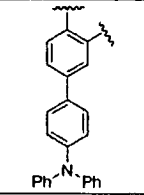
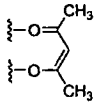

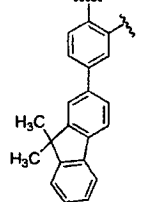
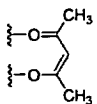

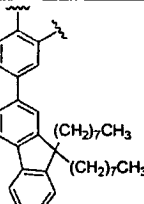
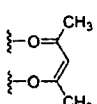

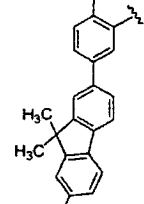
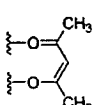

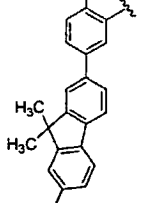
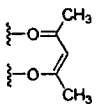

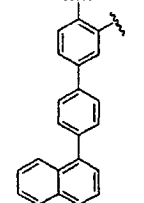
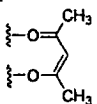
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
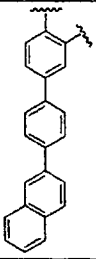
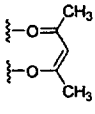

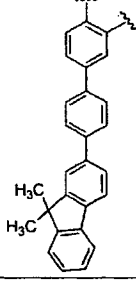
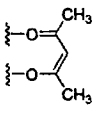

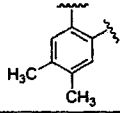
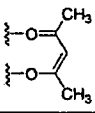

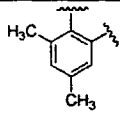
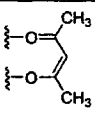

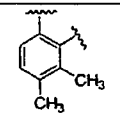
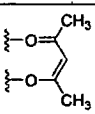

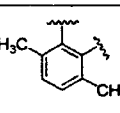
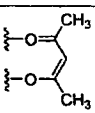

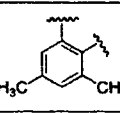
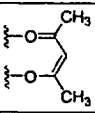

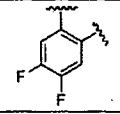
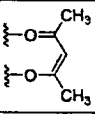

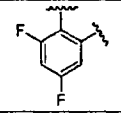
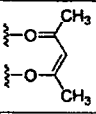
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



































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
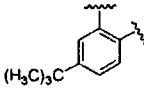
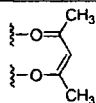


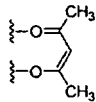

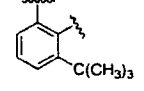
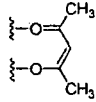

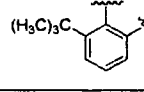
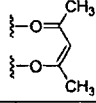

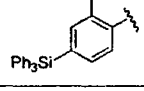
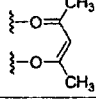

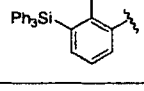
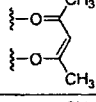

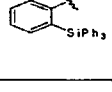
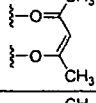

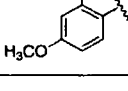
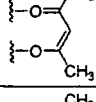

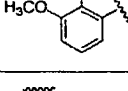
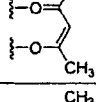

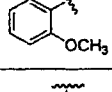
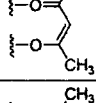

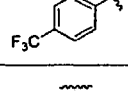
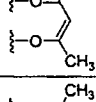

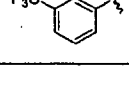
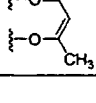
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
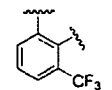


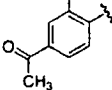


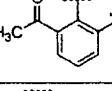


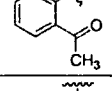


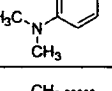


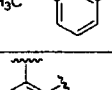


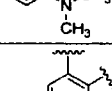


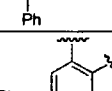


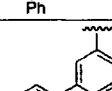


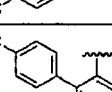


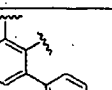


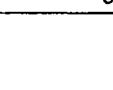

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
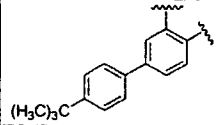
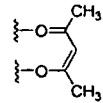

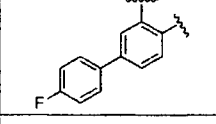
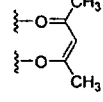

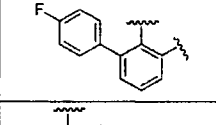
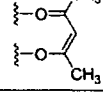

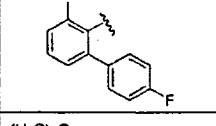
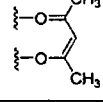

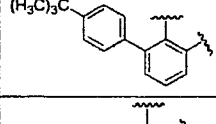
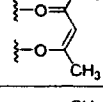

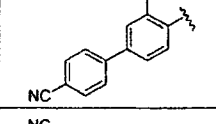
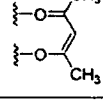

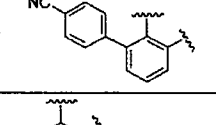
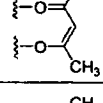

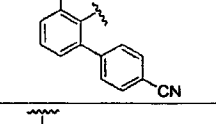
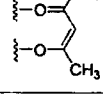

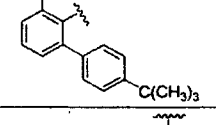
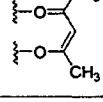

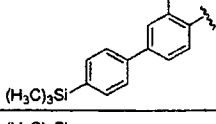
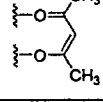

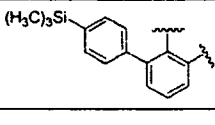
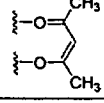
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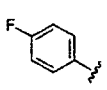
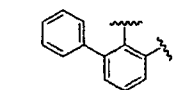
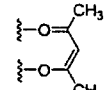
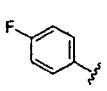
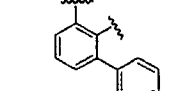
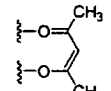
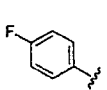
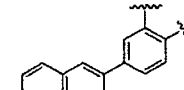
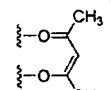
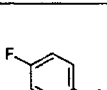
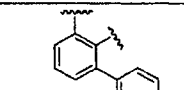
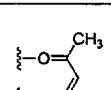
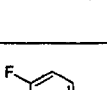
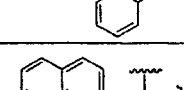
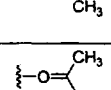
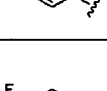
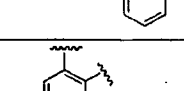
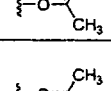
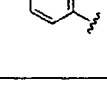
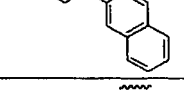
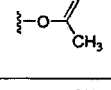
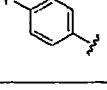
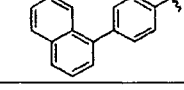
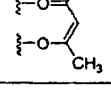
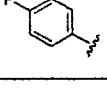
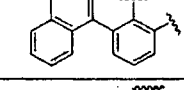
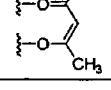
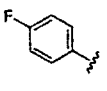
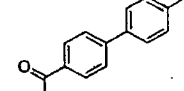
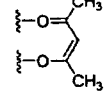
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
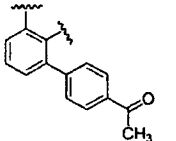
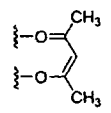

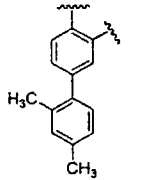
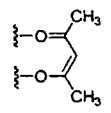

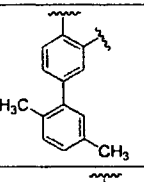
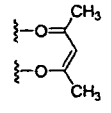

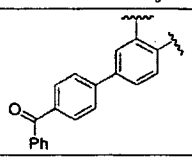
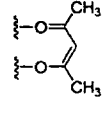

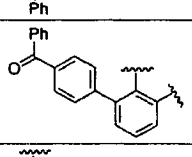
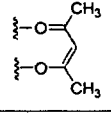

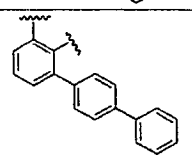
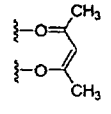

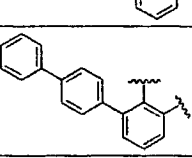
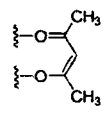

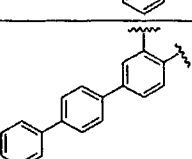
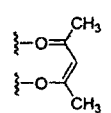

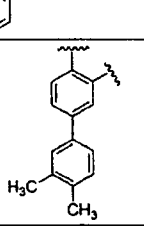
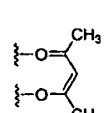
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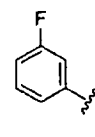
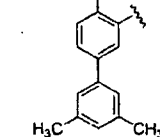
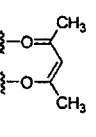
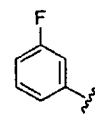
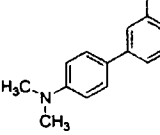
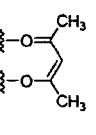
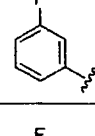
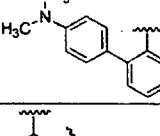
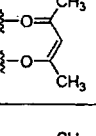
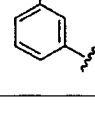
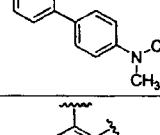
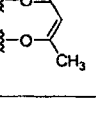
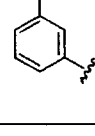
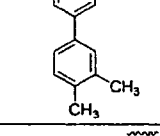
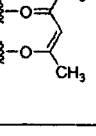
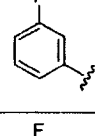
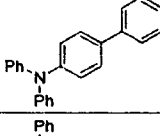
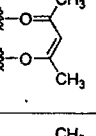
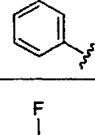
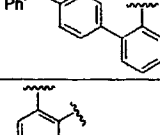
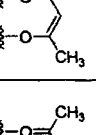
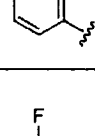
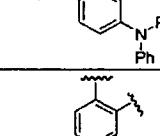
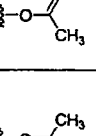
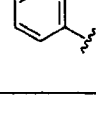
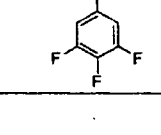
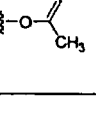
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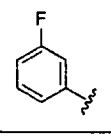
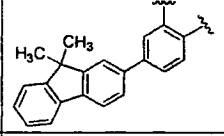
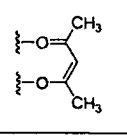
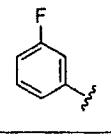
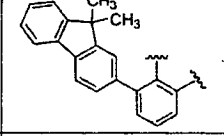
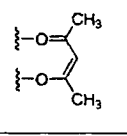
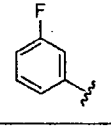
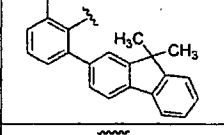
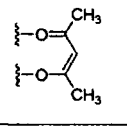
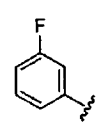
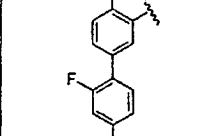
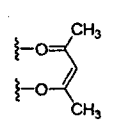
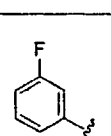
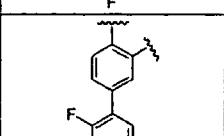
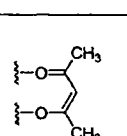
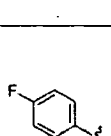
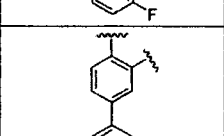
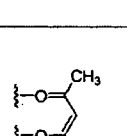
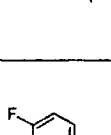
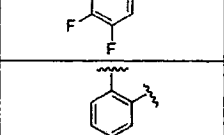
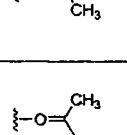
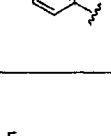
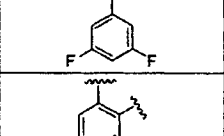
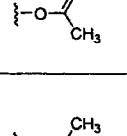
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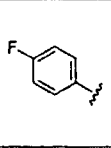
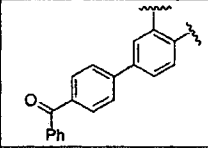
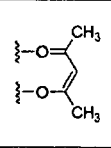
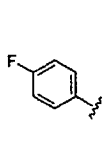
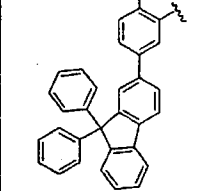
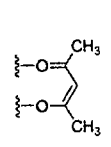
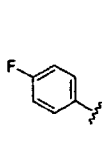
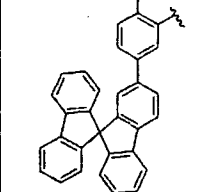
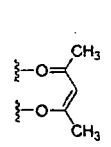
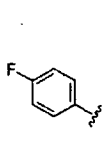
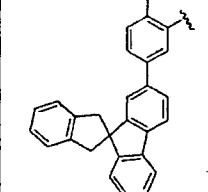
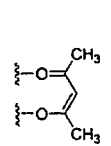
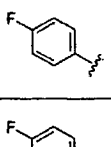
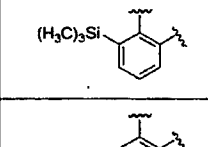
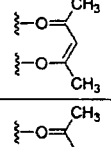
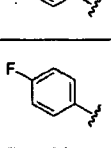
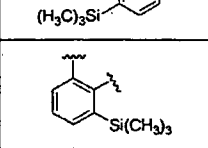
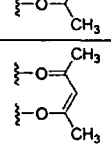
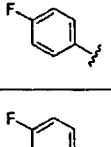
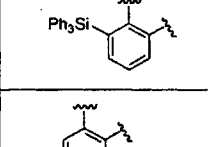
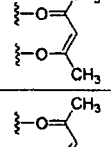
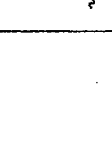
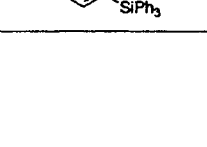
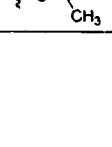



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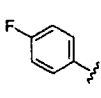
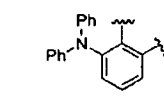

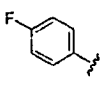
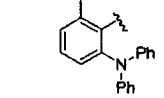

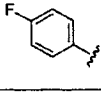
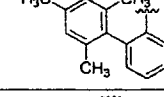

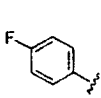
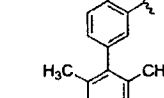

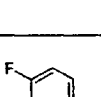
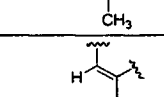

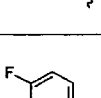
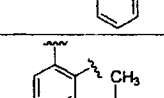

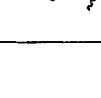
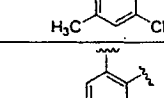

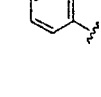
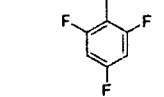

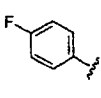
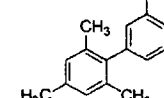

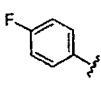
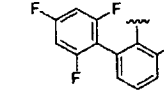

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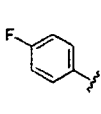
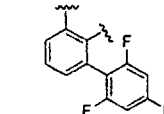
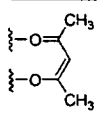
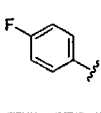
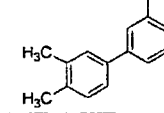
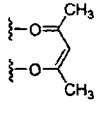
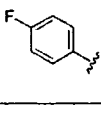
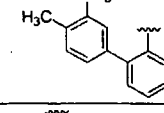
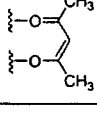
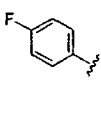
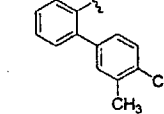
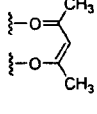
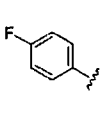
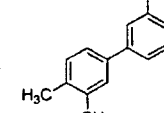
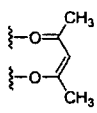
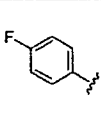
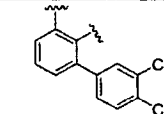
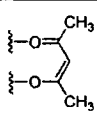
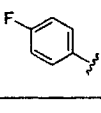
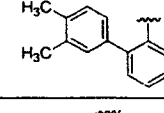
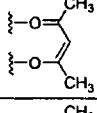
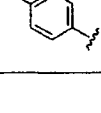
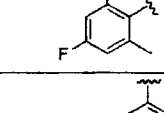
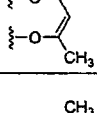
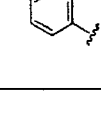
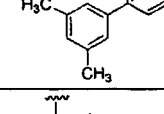
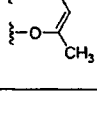
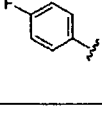
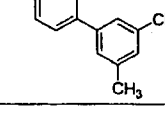
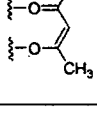
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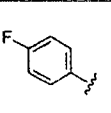
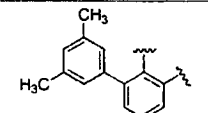
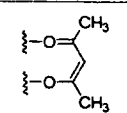
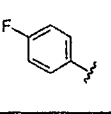
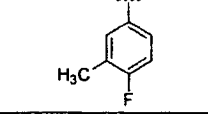
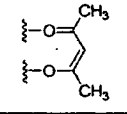
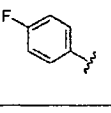
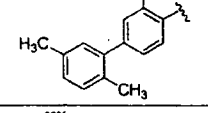
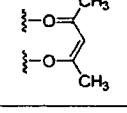
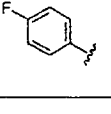
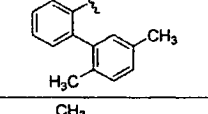
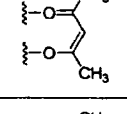
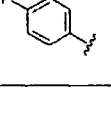
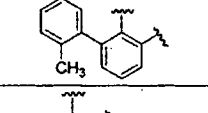
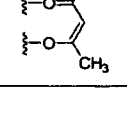
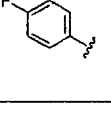
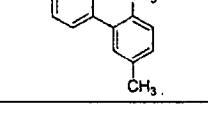
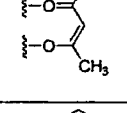
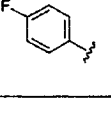
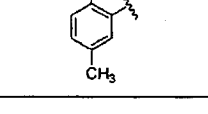
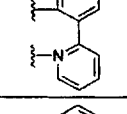
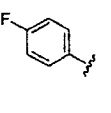
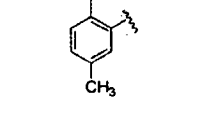
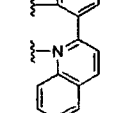
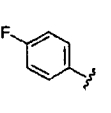
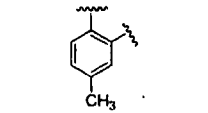
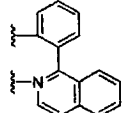
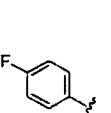
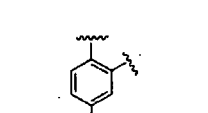
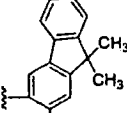
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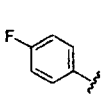
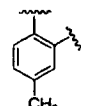
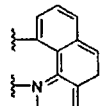
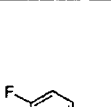
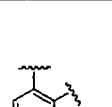
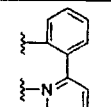

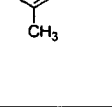
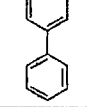
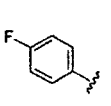
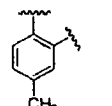
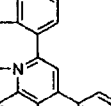
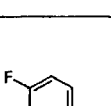
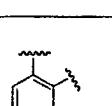
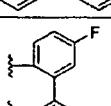

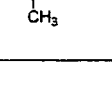
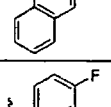
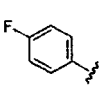
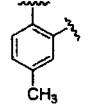
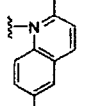
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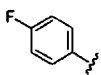
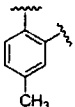
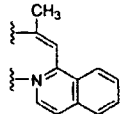
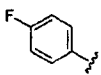
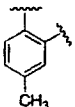
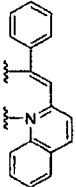
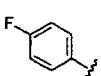
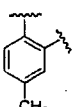
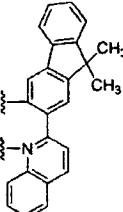
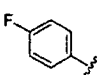
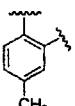
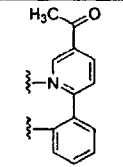
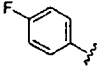
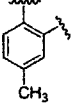
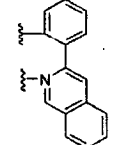
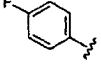
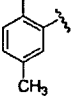
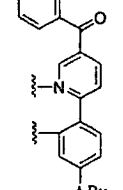
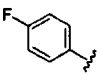
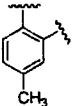
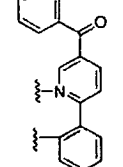
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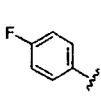
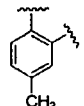
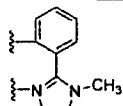
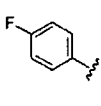
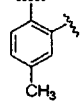
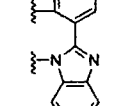
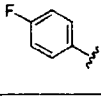
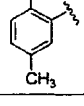
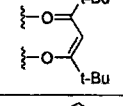
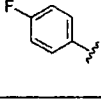
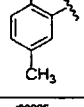
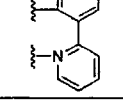
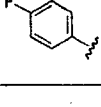
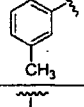
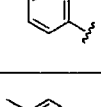
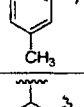
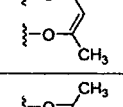
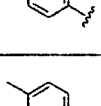
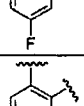
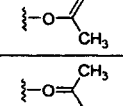
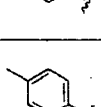
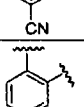
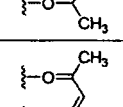
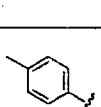
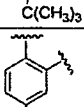
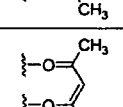
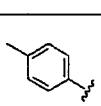
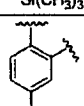
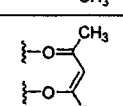



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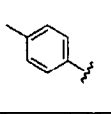
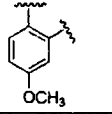
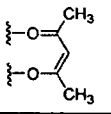
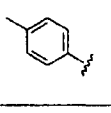
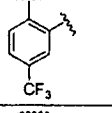
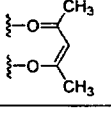
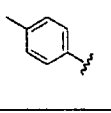
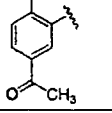
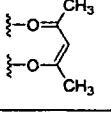
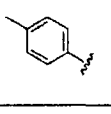
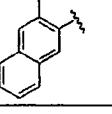
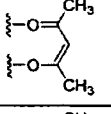
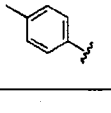
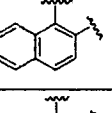
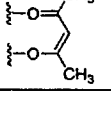
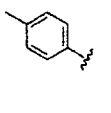
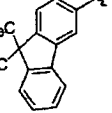
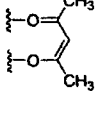
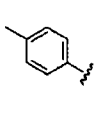
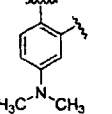
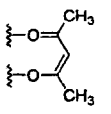
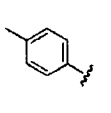
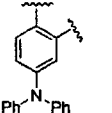
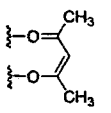
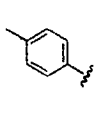
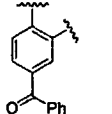
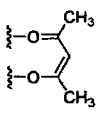
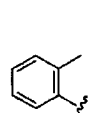
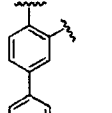
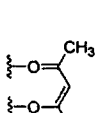
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
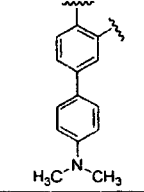
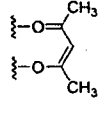

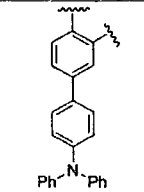
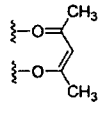

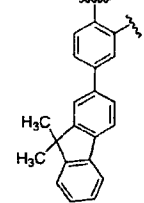
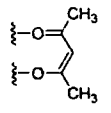

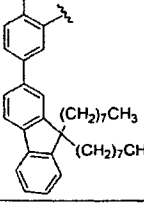
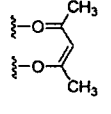

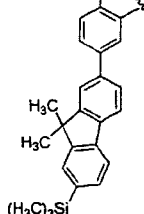
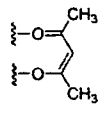

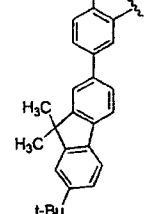
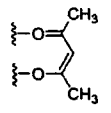
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
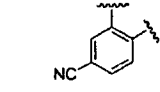


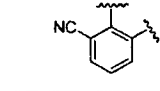


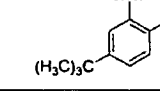


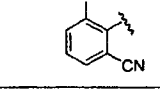


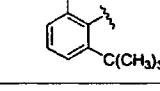


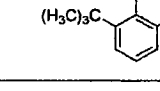


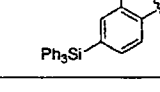


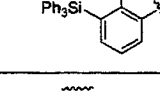


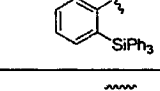


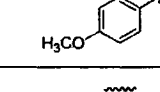


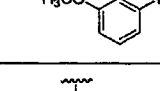


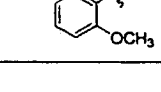

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
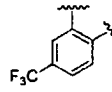
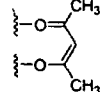

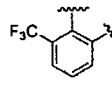
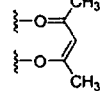

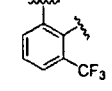
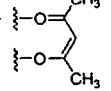

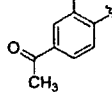
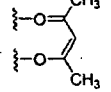

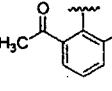
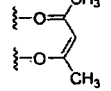

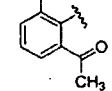
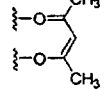

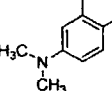
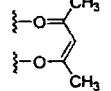

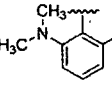
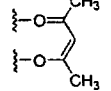

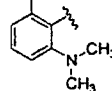
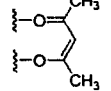

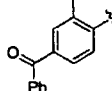
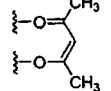

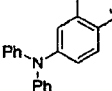
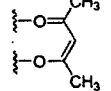

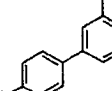
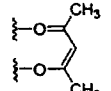
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
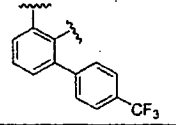
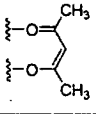

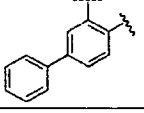
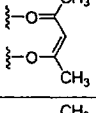

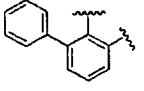
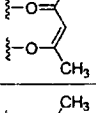

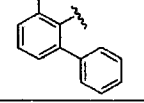
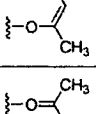

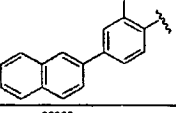
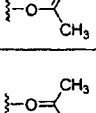

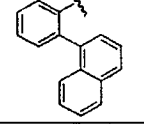
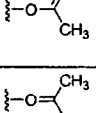

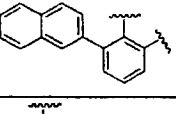
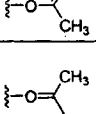

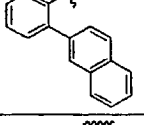
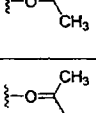

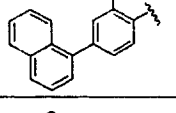
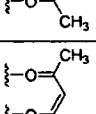

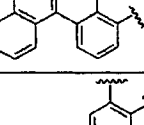
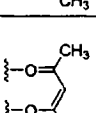

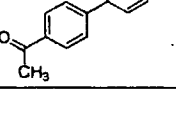
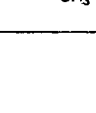
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
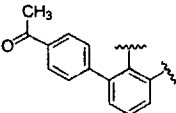
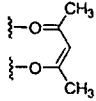

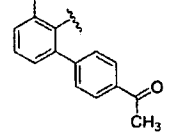
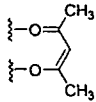

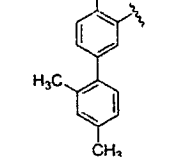
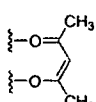

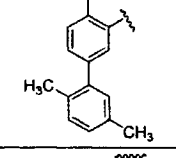
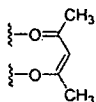

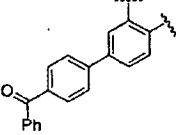
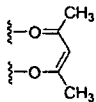

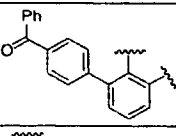
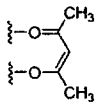

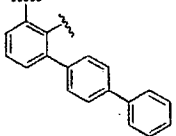
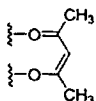

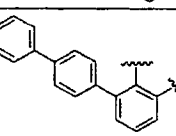
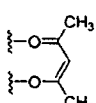

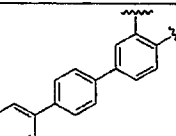
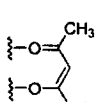
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
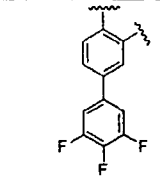
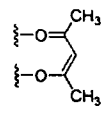

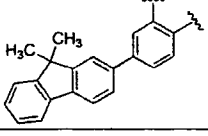
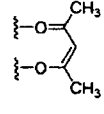

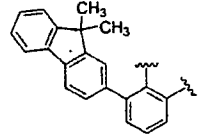
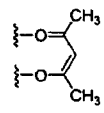

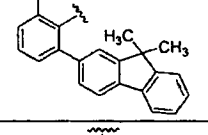
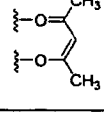

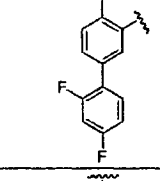
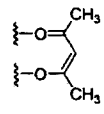

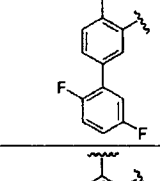
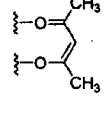

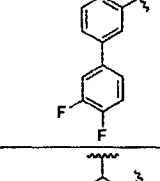
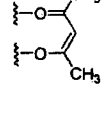

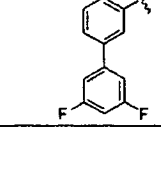
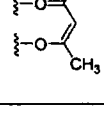
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
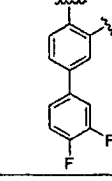
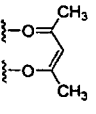

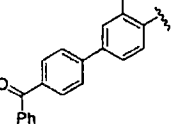
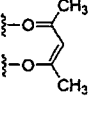

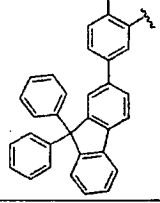
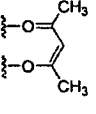

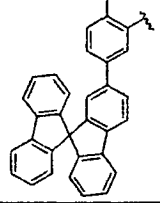
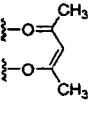

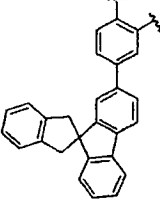
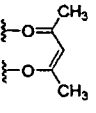

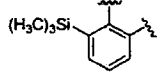
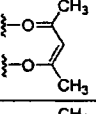

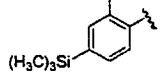
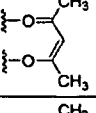

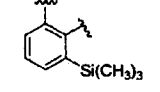
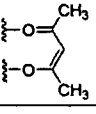
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
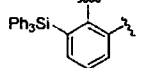
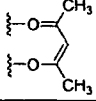

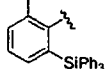
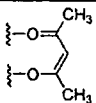

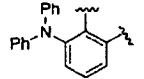
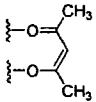

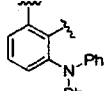
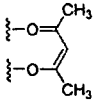

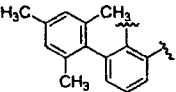
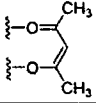

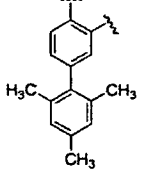
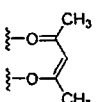

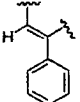
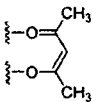

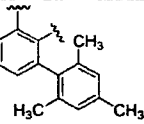
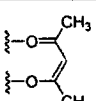

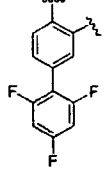
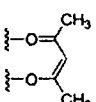

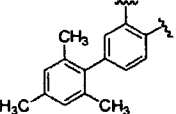
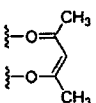

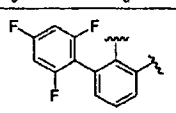
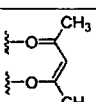
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
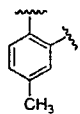
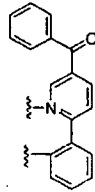

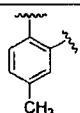
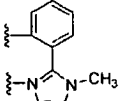

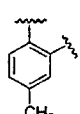
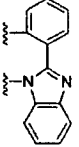

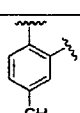
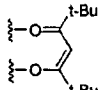

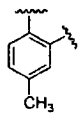
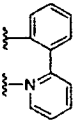

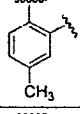

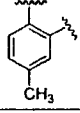
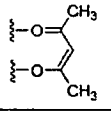

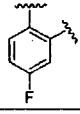
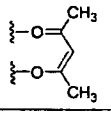

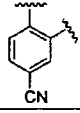
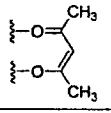

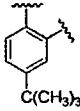
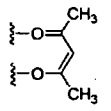
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
































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
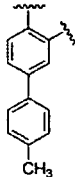
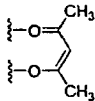

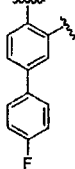


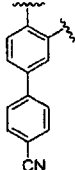
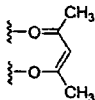

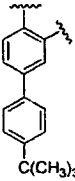
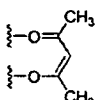

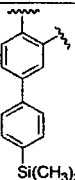


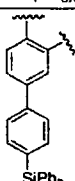


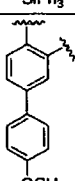

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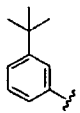
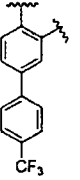
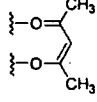
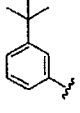
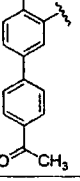
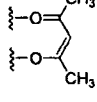
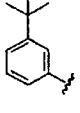
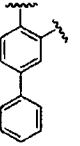
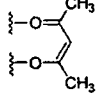
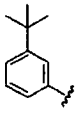
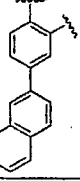
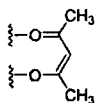
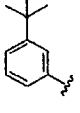
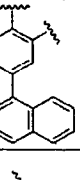
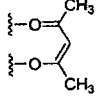
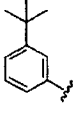
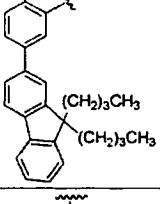
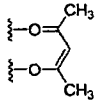
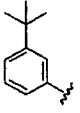
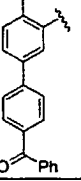
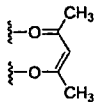
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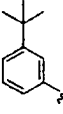
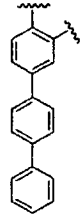
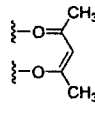
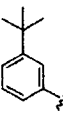
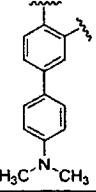
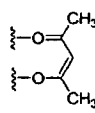
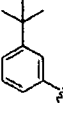
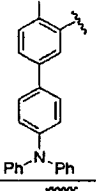
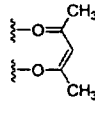
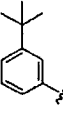
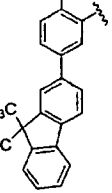
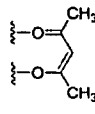
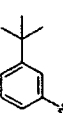
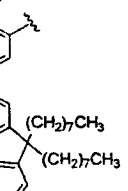
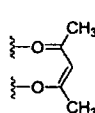

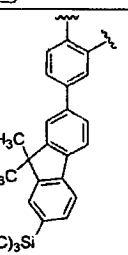
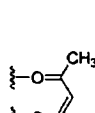
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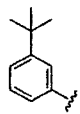
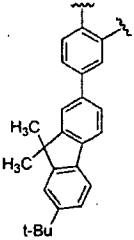
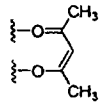
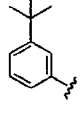
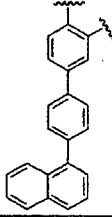
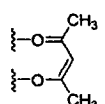
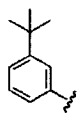
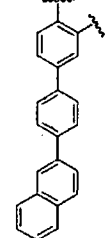
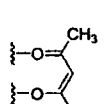
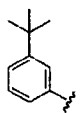
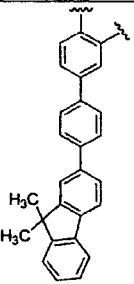
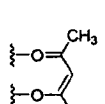
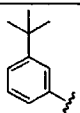
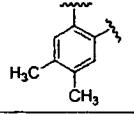
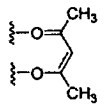
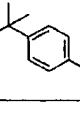
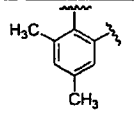
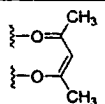
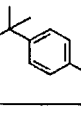
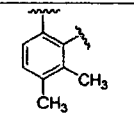
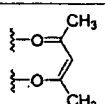
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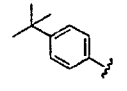
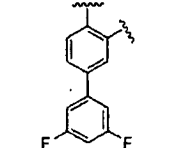
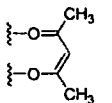
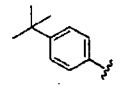
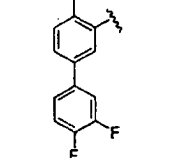
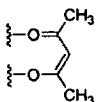
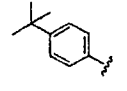
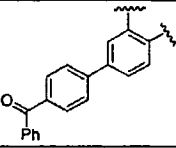
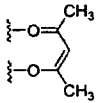
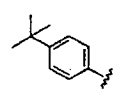
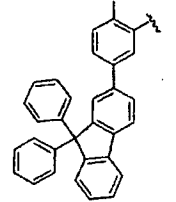
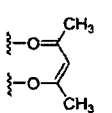
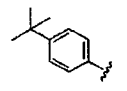
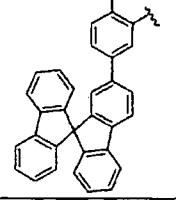
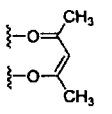
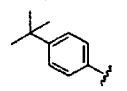
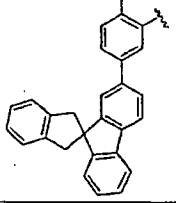
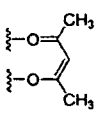
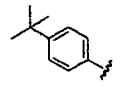
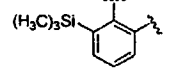
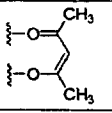
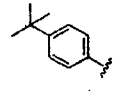
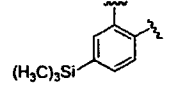
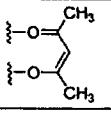
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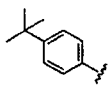
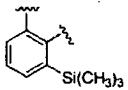
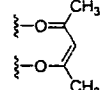
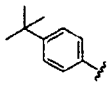
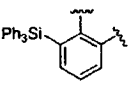
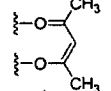
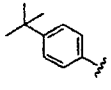
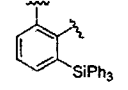
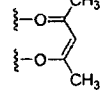
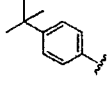
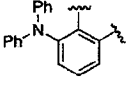
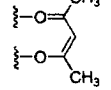
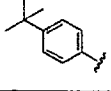
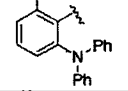
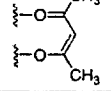
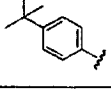
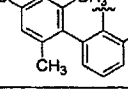
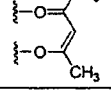
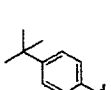
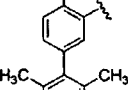
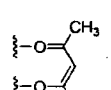
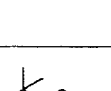
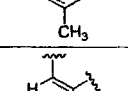
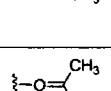
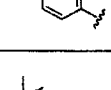
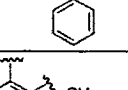
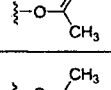
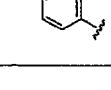
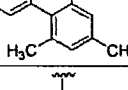
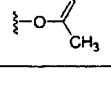
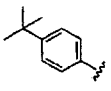
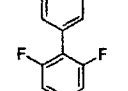
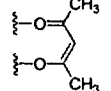
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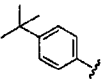
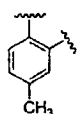
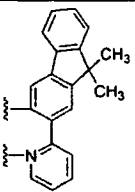
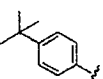
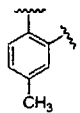
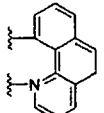
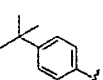
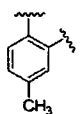
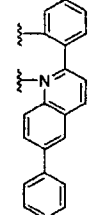
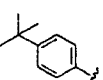
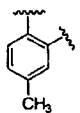
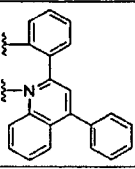
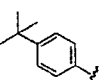
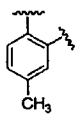
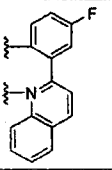
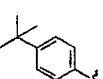
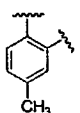
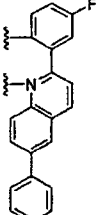
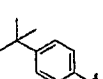
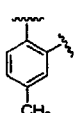
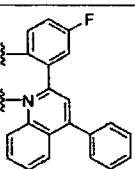
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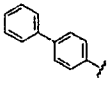

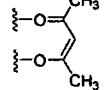
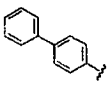

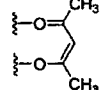
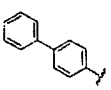

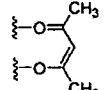
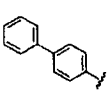

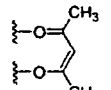
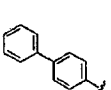

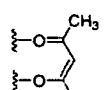
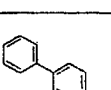

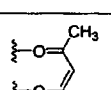
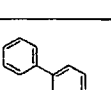

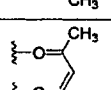
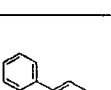

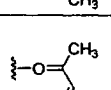
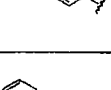

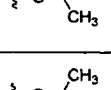
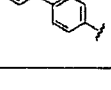

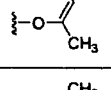
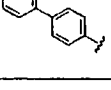

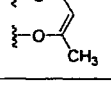
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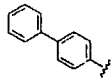
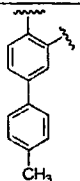
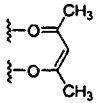
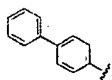
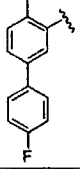
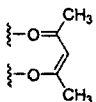
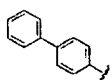
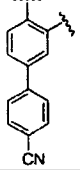
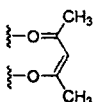
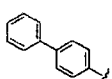
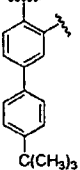
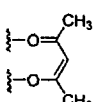
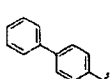
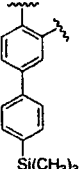
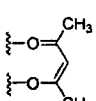
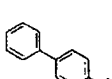
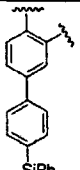
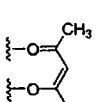
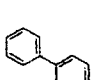
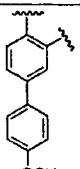
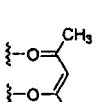
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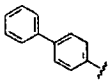
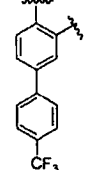
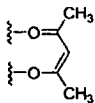
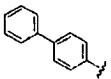
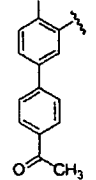
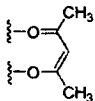
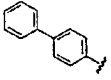
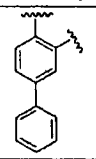
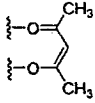
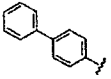
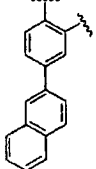
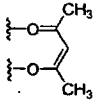
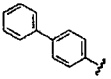
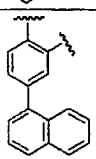
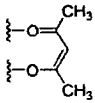
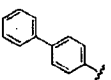
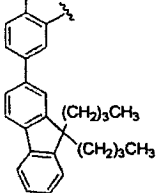
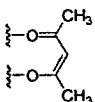
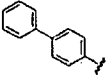
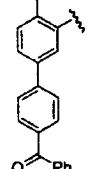
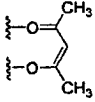
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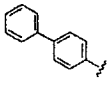
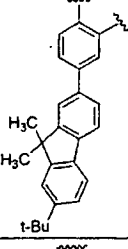
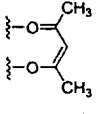
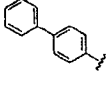
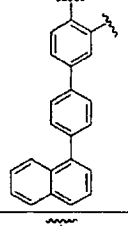
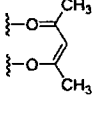
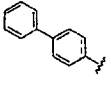
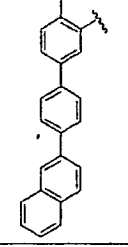
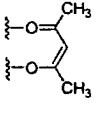
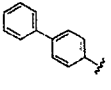
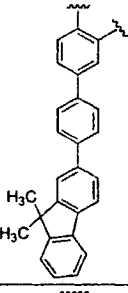
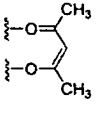
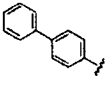
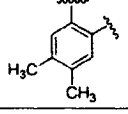
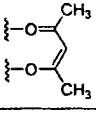
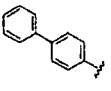
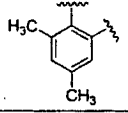
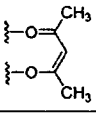
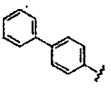
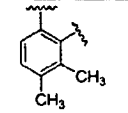
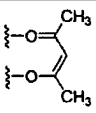
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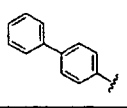
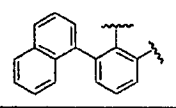
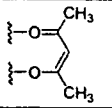
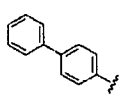
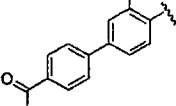
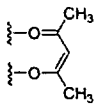
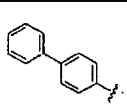
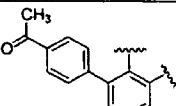
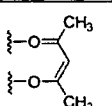
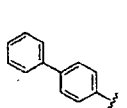
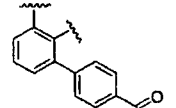
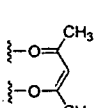
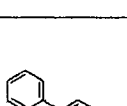
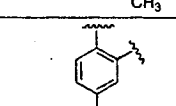
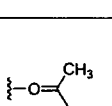
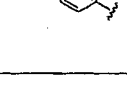
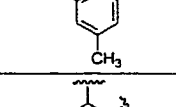
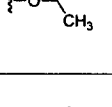
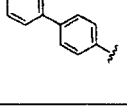
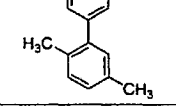
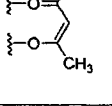
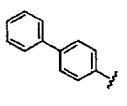
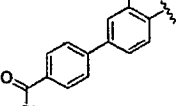
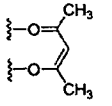
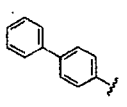
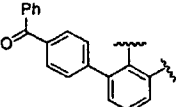
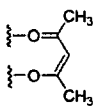
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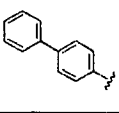
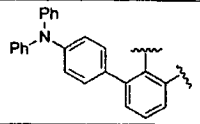
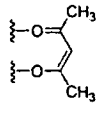
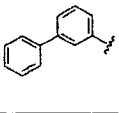
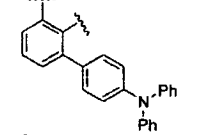
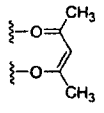
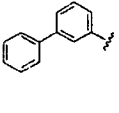
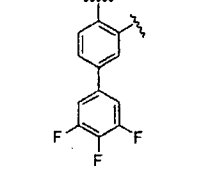
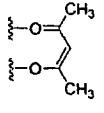
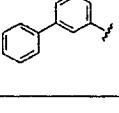
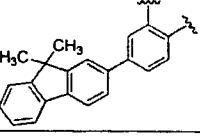
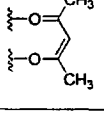
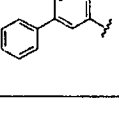
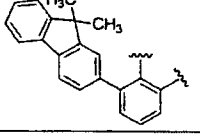
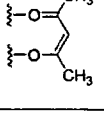
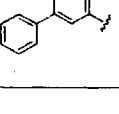
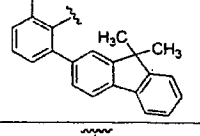
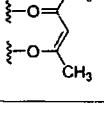
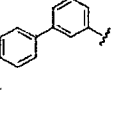
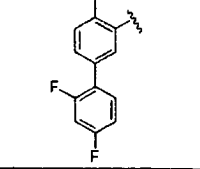
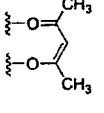
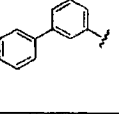
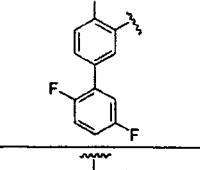
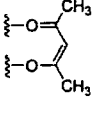
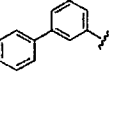
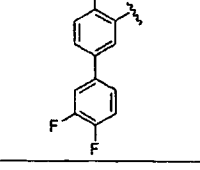
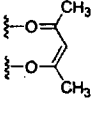
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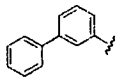
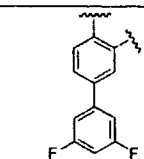
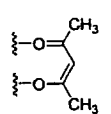
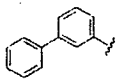
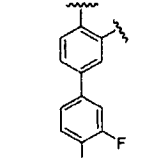
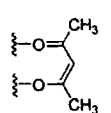
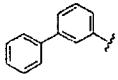
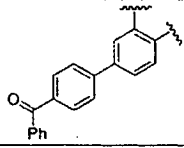
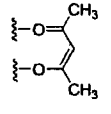
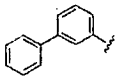
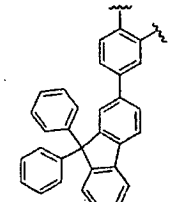
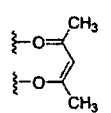
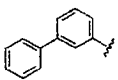
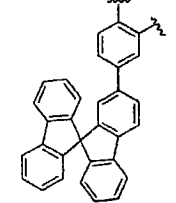
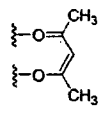
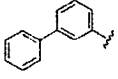
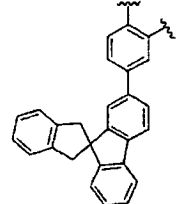
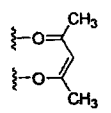
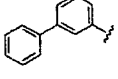
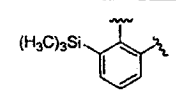
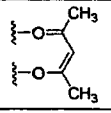
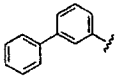
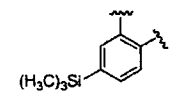
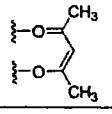
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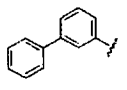
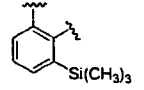
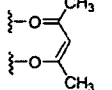
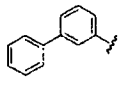
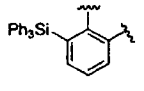
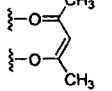
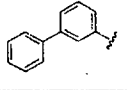
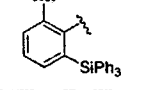
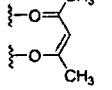
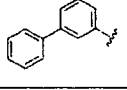
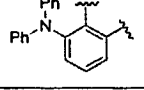
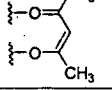
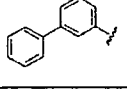
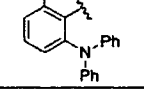
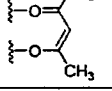
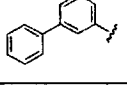
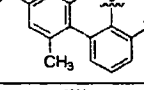
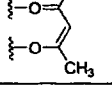
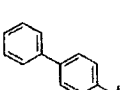
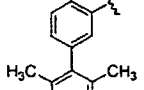
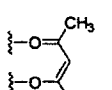
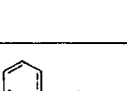
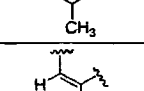
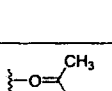
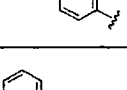
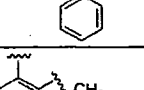
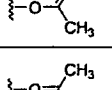
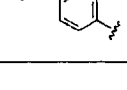
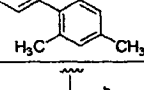
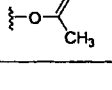
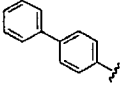
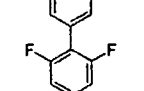
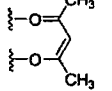
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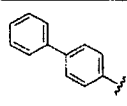
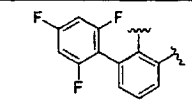
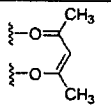
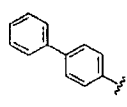
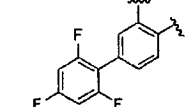
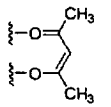
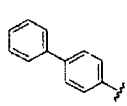
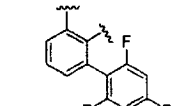
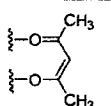
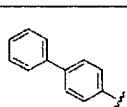
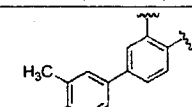
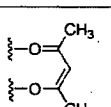
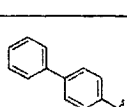
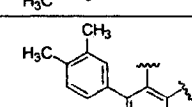
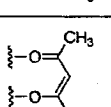
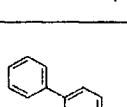
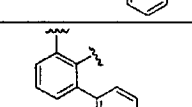
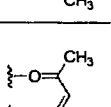
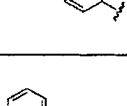
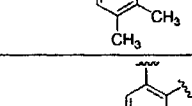
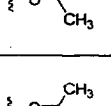
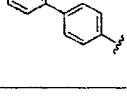
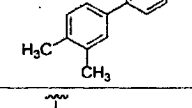
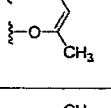
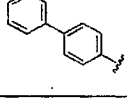
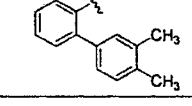
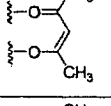
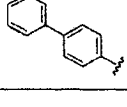
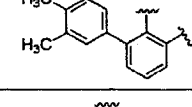
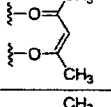
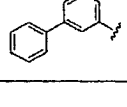
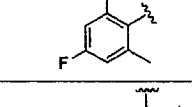
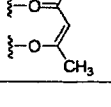
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936	H	H	H	H				2
937	H	H	H	H				2
938	H	H	H	H				2

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939	H	H	H	H				2
940	H	H	H	H				2
941	H	H	H	H				2
942	H	H	H	H				2
943	H	H	H	H				2
944	H	H	H	H				2
945	H	H	H	H				2
946	H	H	H	H				2
947	H	H	H	H				2
948	H	H	H	H				2
949	H	H	H	H				2

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10  
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950	H	H	H	H				2
951	H	H	H	H				2
952	H	H	H	H				2
953	H	H	H	H				2
954	H	H	H	H				2
955	H	H	H	H				2
956	H	H	H	H				2
957	H	H	H	H				2
958	H	H	H	H				2
959	H	H	H	H				2
960	H	H	H	H				2

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5	961	H	H	H	H				2
10	962	H	H	H	H				2
15	963	H	H	H	H				2
20	964	H	H	H	H				2
25	965	H	H	H	H				2
30	966	H	H	H	H				2
35	967	H	H	H	H				2
40	968	H	H	H	H				2
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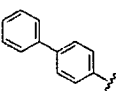
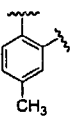
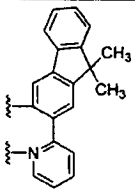
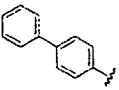
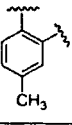
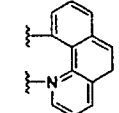
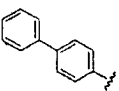
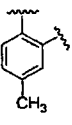
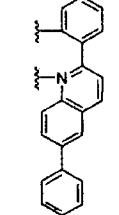
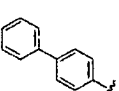
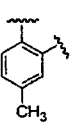
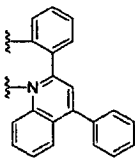
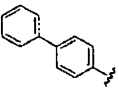
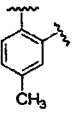
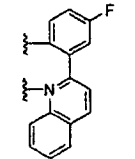
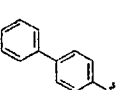
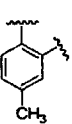
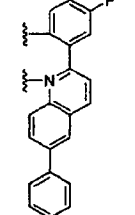
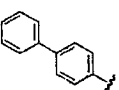
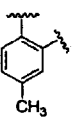
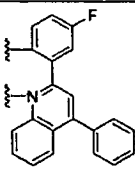
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974	H	H	H	H				2
975	H	H	H	H				2
976	H	H	H	H				2
977	H	H	H	H				2

5	978	H	H	H	H				2
10	979	H	H	H	H				2
15	980	H	H	H	H				2
20	981	H	H	H	H				2
25	982	H	H	H	H				2
30	983	H	H	H	H				2
35	984	H	H	H	H				2
40									
45									
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55									

5	985	H	H	H	H				2
10	986	H	H	H	H				2
15	987	H	H	H	H				2
20	988	H	H	H	H				2
25	989	H	H	H	H				1
30	990	H	H	H	H			-	3
35	991	H	H	H	H				2
40	992	H	H	H	H				2
45	993	H	H	H	H				2
50	994	H	H	H	H				2

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

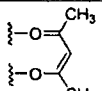


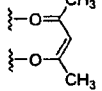


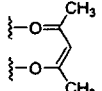


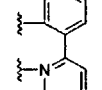




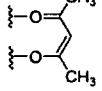


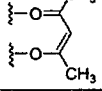


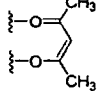


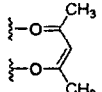


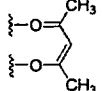
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996	H	H	H	H				2
997	H	H	H	H				2
998	H	H	H	H				2
999	H	H	H	H			-	3
1000	-CH <sub>3</sub>	H	H	H				2
1001	H	-CH <sub>3</sub>	H	H				2
1002	H	H	F	H				2
1003	H	H	H	H				2
1004	-CH <sub>3</sub>	H	H	H				2

Table 2

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Comp. No.	<sup>1</sup> H NMR(CDCl <sub>3</sub> , 200 MHz)	MS/FAR	
		found	calculated
1	δ = 8.07(m, 2H), 7.87(m, 2H), 7.65-7.57(m, 4H). 7.48-7.42(m, 6H). 7.32(m, 4H). 7.22-7.15(m, 6H). 4.59(s, 1H). 2.32(s, 6H), 2.35(s, 6H), 2.09(s, 6H)	908	908.3
4	δ = 8.07(m, 2H), 7.01(m, 2H), 7.65-7.57(m, 4H). 7.48-7.32(m, 14H), 7.22(m, 2H), 4.60(s, 1H), 2.33(s, 6H). 2.09(s, 6H), 1.34(s, 18H)	992	992.39
11	δ = 8.06(m, 2H). 7.69-7.57(m, 10H). 7.48-7.32(m, 16H), 7.22(m, 2H), 4.58(s, 1H), 2.32(s, 6H), 2.10(s, 6H)	980	980.3
33	δ = 8.81(m, 2H), 8.03-7.87(m, 12H), 7.77-7.74(m, 4H), 7.63-7.41(m, 17H). 7.38(m, 2H), 7.28(m, 2H), 2.31(s, 6H), 2.09(s, 6H), 1.72(s, 12H)	1264	1264.45
43	δ = 8.06(m, 2H), 7.65(m, 2H), 7.57(m, 2H). 7.48-7.42(m, 6H), 7.32(m, 4H), 7.22(m, 2H), 7.03(m, 2H), 6.96(m, 2H), 4.59(s, 1H), 2.35(s, 12H). 2.32(s, 6H), 2.09(s, 6H)	936	936.33
50	δ = 8.06(m, 2H), 7.65-7.57(m, 4H), 7.48-7.42(m, 6H). 7.32(m, 4H), 7.22(m, 2H), 7.04-6.97(m, 4H), 4.58(s, 1H), 2.32(s, 6H), 2.08(s, 6H)	952	952.23

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(continued)

Comp. No.	<sup>1</sup> H NMR(CDCl <sub>3</sub> , 200 MHz)	MS/FAR	
		found	calculated
5 69	δ = 8.18(d, 2H), 8.06(m, 2H), 7.65-7.57(m, 4H), 7.48-7.42(m, 8H), 7.32-7.22(m, 8H), 4.60(s, 1H), 2.33(s, 6H), 2.09(s, 6H)	1016	1016.24
10 104	δ = 8.21(m, 2H), 8.03-7.99(m, 4H), 7.74(m, 2H), 7.60-7.41(m, 26H), 4.59(s, 1H), 2.31(s, 6H), 2.09(s, 6H)	1032	1032.33
10 117	δ = 8.06-8.05(m, 4H), 7.65-7.57(m, 8H), 7.48-7.42(m, 6H), 7.32(m, 4H), 7.22-7.16(m, 4H), 7.00(m, 2H), 6.90(m, 2H), 4.58(s, 1H), 2.32(s, 6H), 2.35(s, 12H), 2.09(s, 6H)	1088	1088.39
15 177	δ = 8.30(m, 1H), 8.18(m, 2H), 8.10-7.98(m, 7H), 7.78-7.74(m, 3H), 7.60-7.32(m, 21H), 2.34(s, 6H), 2.31(s, 6H)	1013	1013.33
15 184	δ = 8.21-8.18(m, 3H), 8.10-7.99(m, 7H), 7.90(m, 1H), 7.74(m, 2H), 7.59-7.51(m, 15H), 7.41-7.35(m, 8H), 7.26(m, 1H), 2.34(s, 6H), 2.31(s, 6H)	1107	1107.35
20 192	δ = 8.64(m, 1H), 8.34(m, 1H), 8.22-8.18(m, 3H), 8.03-7.99(m, 4H), 7.89-7.87(m, 3H), 7.74-7.51(m, 15H), 7.41(m, 4H), 7.32(m, 4H), 2.34(s, 6H), 2.31(s, 6H), 1.35(s, 9H)	1123	1123.41
20 198	δ = 8.18(m, 3H), 8.03-7.99(m, 6H), 7.74(m, 3H), 7.59(m, 18H), 7.32(m, 6H), 2.34(s, 9H), 2.31(s, 9H)	1117	1117.39
25 241	δ = 8.06(m, 2H), 7.65(m, 2H), 7.57(m, 2H), 7.46-7.42(m, 6H), 7.03-7.96(m, 8H), 4.58(s, 1H), 2.35(s, 12H), 2.32(s, 6H), 2.09(s, 6H)	972	972.31
25 244	δ = 8.03-7.99(m, 4H), 7.74-7.71(m, 6H), 7.59(m, 2H), 7.49(m, 2H), 7.28(m, 2H), 7.10(m, 2H), 6.77(m, 2H), 4.60(s, 1H), 2.31(s, 6H), 2.09(s, 6H)	988	988.21
30 309	δ = 8.21(m, 2H), 8.06(m, 2H), 7.67-7.32(m, 28H), 7.03(m, 4H), 4.58(s, 1H), 2.33(s, 6H), 2.09(s, 6H)	1168	1168
30 406	δ = 8.85(m, 2H), 8.03-7.95(m, 10H), 7.74(m, 2H), 7.59(m, 6H), 7.33-7.29(m, 8H), 4.59(s, 1H), 2.34(s, 6H), 2.31(s, 6H), 2.09(s, 6H)	1008	1008.33
35 437	δ = 8.06-7.99(m, 4H), 7.79-7.74(m, 4H), 7.59(m, 2H), 7.39(m, 4H), 7.19-7.13(m, 6H), 4.59(s, 1H), 2.59(s, 6H), 2.34(s, 12H), 2.31(s, 6H), 2.09(s, 6H)	964	964.36
35 620	δ = 8.81(m, 2H), 8.03-7.92(m, 16H), 7.74-7.73(m, 4H), 7.59-7.58(m, 8H), 7.43-7.42(m, 4H), 7.32(m, 2H), 4.59(s, 1H), 2.31(s, 6H), 2.09(s, 6H), 1.35(s, 18H)	1244	1244.48
40 648	δ = 8.30(m, 2H), 8.03-7.99(m, 4H), 7.74(m, 2H), 7.59(m, 2H), 7.42-7.37(m, 10H), 7.25(m, 2H), 4.59(s, 1H), 2.34(s, 6H), 2.31(s, 6H), 2.09(s, 6H), 1.35(s, 18H)	1020	1020.42
40 766	δ = 8.21(m, 2H), 8.06(m, 2H), 7.65-7.50(m, 6H), 7.42-7.35(m, 12H), 7.16(s, 2H), 7.00(m, 2H), 6.90(m, 2H), 4.62(s, 1H), 2.37(s, 12H), 2.34(s, 6H), 2.08(s, 6H), 1.35(s, 18H)	1200	1200.51
45 838	δ = 8.03-7.99(m, 4H), 7.74(m, 2H), 7.59-7.41(m, 12H), 7.25(m, 8H), 7.10(m, 2H), 6.77(m, 2H), 4.59(s, 1H), 2.31(s, 6H), 2.09(s, 6H)	1104	1104.29
45 842	δ = 8.03-7.99(m, 4H), 7.74(m, 2H), 7.59-7.41(m, 12H), 7.31-7.21(m, 12H), 4.59(s, 1H), 2.31(s, 6H), 2.09(s, 6H)	1104	1104.29
50 938	δ = 8.21(m, 2H), 8.03-7.99(m, 4H), 7.74-7.70(m, 4H), 7.59-7.41(m, 22H), 4.59(s, 1H), 2.31(s, 6H), 2.09(s, 6H), 0.25(s, 18H)	1176	1176.41
50 993	δ = 8.03-7.94(m, 6H), 7.74(m, 2H), 7.59-7.41(m, 18H), 4.59(s, 1H), 2.31(s, 6H), 2.09(s, 6H)	880	880.26
55 994	δ = 8.03-7.94(m, 6H), 7.74(m, 2H), 7.59-7.47(m, 8H), 7.39-7.30(m, 8H), 4.59(s, 1H), 2.31(s, 6H), 2.09(s, 6H)	916	916.25

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(continued)

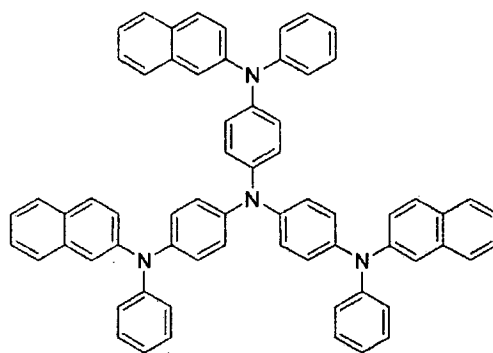
Comp. No.	<sup>1</sup> H NMR(CDCl <sub>3</sub> , 200 MHz)	MS/FAR	
		found	calculated
999	δ = 8.03-7.99(m, 9H), 7.74(m, 3H), 7.59-7.47(m, 27H), 2.31(s, 9H)	1075	1075.35

[Example 1] Manufacture of an OLED (1)

**[0061]** An OLED device was manufactured by using a red phosphorescent compound according to the invention.

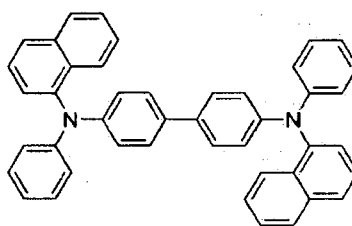
**[0062]** First, a transparent electrode ITO thin film (15 Ω/□) (2) prepared from glass for OLED (produced by Samsung Corning) (1) was subjected to ultrasonic washing with trichloroethylene, acetone, ethanol and distilled water, sequentially, and stored in isopropanol before use.

**[0063]** Then, an ITO substrate was equipped in a substrate folder of a vacuum vapor-deposit device, and 4,4',4"-tris (N,N-(2-naphthyl)-phenylamino)triphenylamine (2-TNATA) was placed in a cell of the vacuum vapor-deposit device, which was then ventilated up to 10<sup>-6</sup> torr of vacuum in the chamber. Electric current was applied to the cell to evaporate 2-TNATA, thereby providing vapor-deposit of a hole injection layer (3) having 60 nm of thickness on the ITO substrate.



2-TNATA

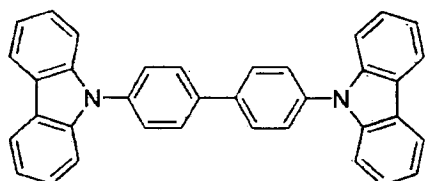
**[0064]** Then, to another cell of the vacuum vapor-deposit device, charged was N,N'-bis(α-naphthyl)-N,N'-diphenyl-4,4'-diamine (NPB), and electric current was applied to the cell to evaporate NPB, thereby providing vapor-deposit of a hole transport layer (4) of 20 nm of thickness on the hole injection layer.



NPB

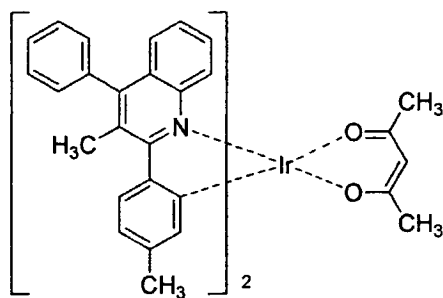
**[0065]** To another cell of said vacuum vapor-deposit device, charged was 4,4'-N,N'-dicarbazole-biphenyl (CBP) as an electroluminescent host material, and an organic electroluminescent compound (Compound 1) according to the present invention was charged to still another cell. The two materials were evaporated at different rates to carry out doping to vapor-deposit an electroluminescent layer (5) having 30 nm of thickness on the hole transport layer. The suitable doping concentration is 4 to 10 mol% on the basis of CBP.

5



CBP

10

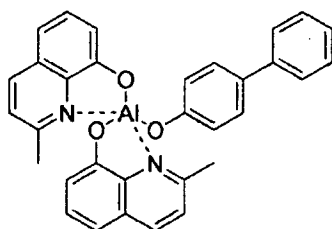


Compound 1

15 **[0066]** Then, on the electroluminescent layer, bis(2-methyl-8-quinolato)(p-phenylphenolato)aluminum (III) (BAIq) was vapor-deposited as a hole blocking layer with a thickness of 10 nm in the same manner for NPB, tris(8-hydroxyquinoline)aluminum (III) (Alq) was vapor-deposited as an electron transport layer (6) with a thickness of 20 nm, and then lithium quinolate (Liq) was vapor-deposited as an electron injection layer (7) with a thickness of 1 to 2 nm. Thereafter, an Al cathode (8) was vapor-deposited with a thickness of 150 nm by using another vacuum vapor-deposit device to manufacture an OLED.

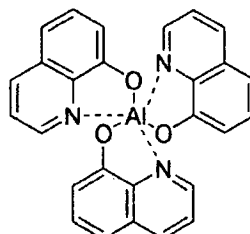
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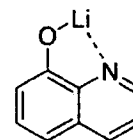


BAIq

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Alq



Liq

35 [Example 2] Manufacture of an OLED (2)

**[0067]** A hole injection layer and a hole transport layer were formed according to the procedure of Example 1, and an electroluminescent layer was vapor-deposited thereon as follows. To another cell of said vacuum vapor-deposit device, charged was H-2 as an electroluminescent host material according to the invention, and an organic electroluminescent compound (Compound 177) according to the present invention was charged to still another cell. The two materials were evaporated at different rates to carry out doping to vapor-deposit an electroluminescent layer (5) having 30 nm of thickness on the hole transport layer. The suitable doping concentration is 4 to 10 mol% on the basis of the host. Then, a hole blocking layer, an electron transport layer and an electron injection layer were vapor-deposited according to the same procedure as in Example 1, and then Al cathode was vapor-deposited in a thickness of 150 nm by using another vacuum vapor-deposit device to manufacture an OLED.

40

45

[Example 3] Manufacture of an OLED (3)

**[0068]** A hole injection layer, a hole transport layer and an electroluminescent layer were formed according to the same procedure as in Example 2, and then an electron transport layer and an electron injection layer were vapor-deposited. Thereafter, Al cathode was vapor-deposited in a thickness of 150 nm by using another vacuum vapor-deposit device to manufacture an OLED.

50

**[0069]** In order to confirm the performance of the OLED's prepared according to Examples 1 to 3, the luminous efficiencies of the OLED's were measured at 10 mA/cm<sup>2</sup>. Various properties are shown in Tables 3.

55

Table 3

	Material	Host	Hole blocking layer	EL color	Operation voltage	Max. luminous efficiency (cd/A)
5	Compound 1	CBP	BAIq	Red	7.8	10.1
	Compound 44	CBP	BAIq	Red	8.0	10.5
	Compound 49	CBP	BAIq	Red	8.1	11.0
10	Compound 177	CBP	BAIq	Red	8.2	10.8
	Compound 241	CBP	BAIq	Red	8.2	9.4
	Compound 303	CBP	BAIq	Red	7.9	8.8
15	Compound 340	CBP	BAIq	Red	8.0	9.6
	Compound 374	CBP	BAIq	Red	7.9	10.4
	Compound 379	CBP	BAIq	Red	8.0	9.2
	Compound 432	CBP	BAIq	Red	7.8	10.0
20	Compound 578	CBP	BAIq	Red	8.2	9.8
	Compound 584	CBP	BAIq	Red	7.9	11.1
	Compound 589	CBP	BAIq	Red	8.0	11.3
25	Compound 645	CBP	BAIq	Red	8.2	9.4
	Compound 780	CBP	BAIq	Red	7.9	10.2
	Compound 790	CBP	BAIq	Red	8.3	9.3
	Compound 932	CBP	BAIq	Red	8.0	10.0
30	Compound 177	H-2	BAIq	Red	8.0	10.6
	Compound 375	H-7	BAIq	Red	7.8	9.9
	Compound 780	H-40	BAIq	Red	7.8	10.0
35	Compound 49	H-8	-	Red	7.0	10.8
	Compound 303	-12	-	Red	7.4	8.9
	Compound 780	-40	-	Red	7.1	10.2
	Compound 790	-64	-	Red	7.0	9.6

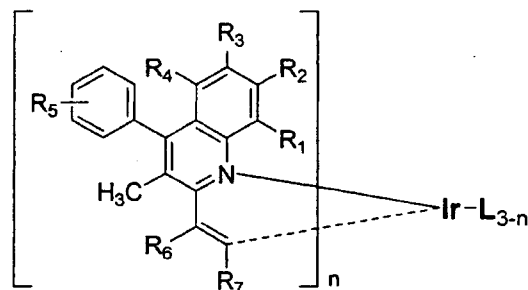
40 **[0070]** For Compound (44), incorporation of dimethyl provided the effect of enhanced efficiency and color coordinate. For Compound (49), incorporation of F to the ligand affected the HOMO level, and provided lowered color coordinate but enhanced luminous efficiency as compared to the compound before incorporating F. Compounds (177), (584) and (589), wherein 2-phenylquinoline, 2-styrylquinoline and phenyl(6-phenylpyridin-3-yl)methanone were incorporated as a subsidiary ligand, respectively, exhibited the highest efficiencies (10.8 cd/A, 11.1 cd/A and 11.3 cd/A, respectively) among the compounds according to the present invention. The compounds wherein ppy or 1-styrylquinoline is incorporated as a subsidiary ligand showed high efficiency of not less than 10 cd/A. Thus, it is found that the subsidiary ligands such as 2-phenylquinoline, 2-styrylquinoline, phenyl(6-phenylpyridin-3-yl)methanone, ppy and 1-styrylquinoline play an important role to enhance the luminous efficiencies of the compounds.

50 **[0071]** With identical device structure, using the host according to the present invention instead of CBP in an EL device did not provide significant change in efficiency, color coordinate and operation voltage from those of CBP. Thus it is anticipated that those hosts can be employed as a phosphorescent host, when being used with dopants according to the invention, instead of CBP as a conventional electroluminescent host. When the host according to the invention is employed without using a hole blocking layer, the device exhibits comparable or higher luminous efficiency as compared to that using a conventional host, and provides decreased power consumption of the OLED due to lowered operation voltage by at least 0.8 V. If the invention is applied to mass production of OLED's, the time for mass production can be also shortened to give great benefit on the commercialization.

## Claims

1. An organic electroluminescent compound represented by Chemical Formula (1):

Chemical Formula 1



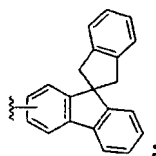
wherein, L is an organic ligand;

$R_1$  through  $R_4$  independently represent hydrogen, (C1-C60)alkyl, (C1-C60)alkoxy, (C3-C60)cycloalkyl, halogen, tri(C1-C60)alkylsilyl or tri(C6-C60)arylsilyl;

$R_5$  represents hydrogen, (C1-C60)alkyl, halogen, or (C6-C60)aryl;

$R_6$  and  $R_7$  independently represent hydrogen, (C1-C60)alkyl with or without halogen substituent(s), (C6-C60)aryl, halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino or di(C6-C60)arylamino, or  $R_6$  and  $R_7$  may be linked via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

the alkyl or aryl of  $R_6$  and  $R_7$ , or the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed therefrom by linkage via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring may be further substituted by one or more substituent(s) selected from (C1-C60)alkyl with or without halogen substituent(s), halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino, di(C6-C60)arylamino, phenyl, naphthyl, anthryl, fluorenyl, spirobifluorenyl and



and

$n$  is an integer from 1 to 3.

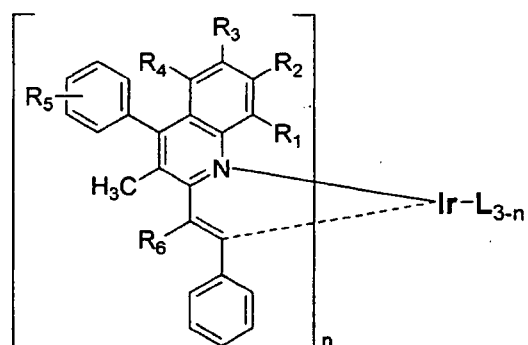
2. The organic electroluminescent compound according to claim 1, which is selected from the compounds represented by one of Chemical Formulas (2) to (7):

Chemical Formula 2

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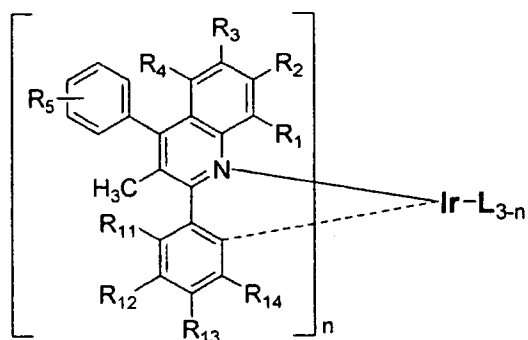


Chemical Formula 3

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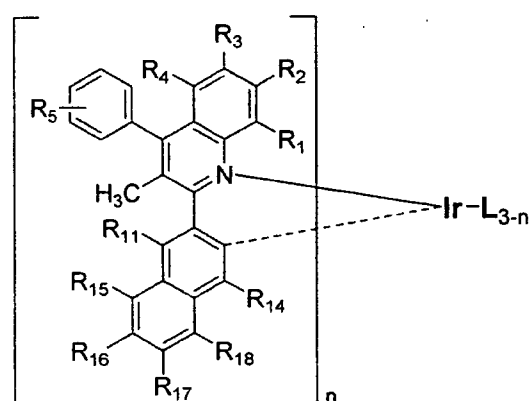
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Chemical Formula 4

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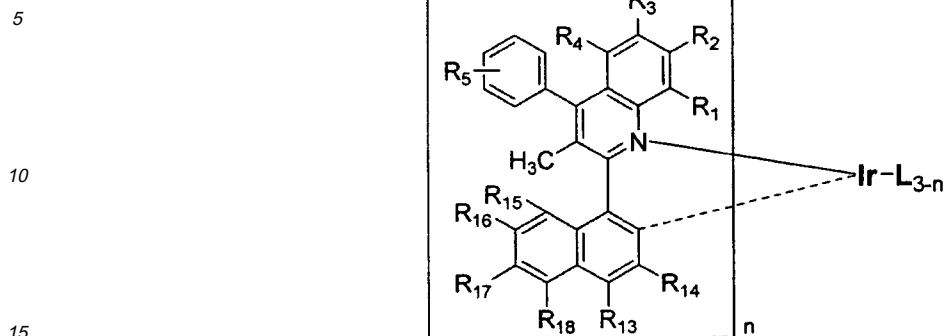
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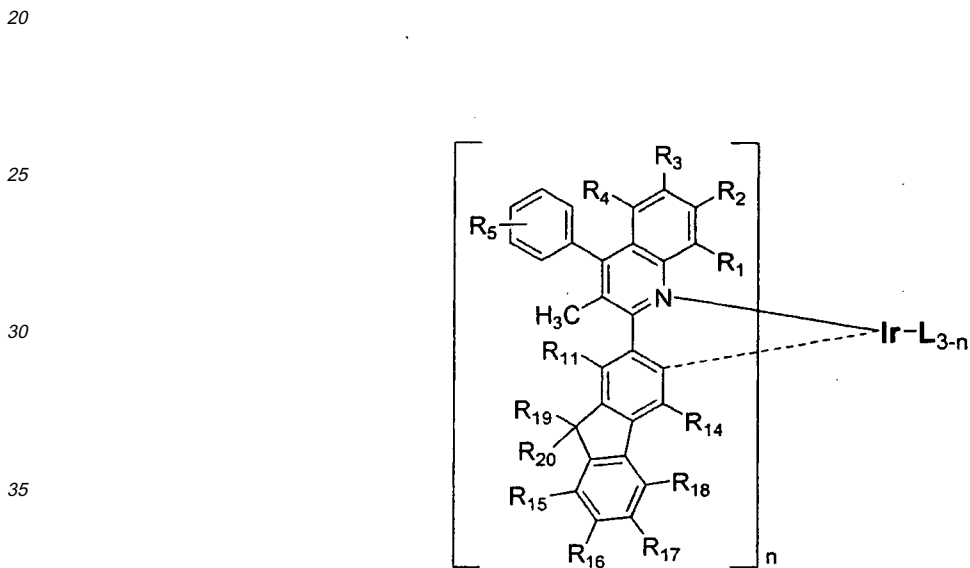


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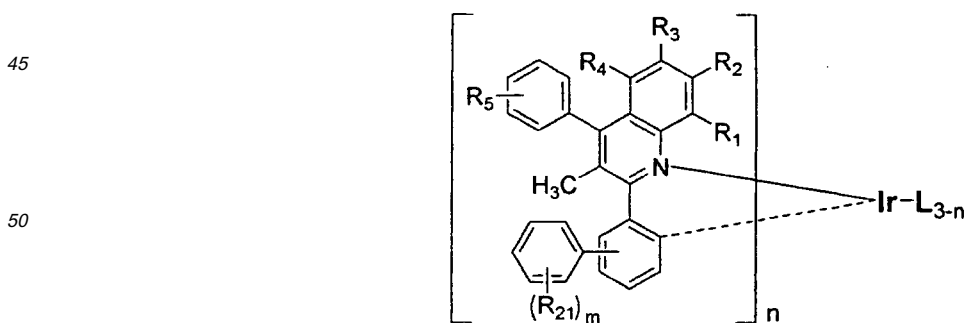
Chemical Formula 5



Chemical Formula 6



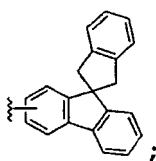
Chemical Formula 7



wherein, L, R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub>, R<sub>5</sub>, R<sub>6</sub> and n are defined as in Chemical Formula (1) of claim 1;  
 R<sub>11</sub> through R<sub>18</sub> independently represent hydrogen, (C1-C60)alkyl, halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino, di(C6-

C60)arylamino, phenyl, naphthyl, anthryl, fluorenyl, spirobifluorenyl or

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the alkyl, phenyl, naphthyl, anthryl or fluorenyl of R<sub>11</sub> through R<sub>18</sub> may be further substituted by one or more substituent(s) selected from (C1-C60)alkyl with or without halogen substituent(s), (C1-C60)alkoxy, halogen, tri (C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino, di (C6-C60)arylamino and (C6-C60)aryl;

15

R<sub>19</sub> and R<sub>20</sub> independently represent hydrogen, (C1-C60)alkyl or (C6-C60)aryl, or R<sub>19</sub> and R<sub>20</sub> may be linked via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring to form an alicyclic ring, a monocyclic or polycyclic aromatic ring;

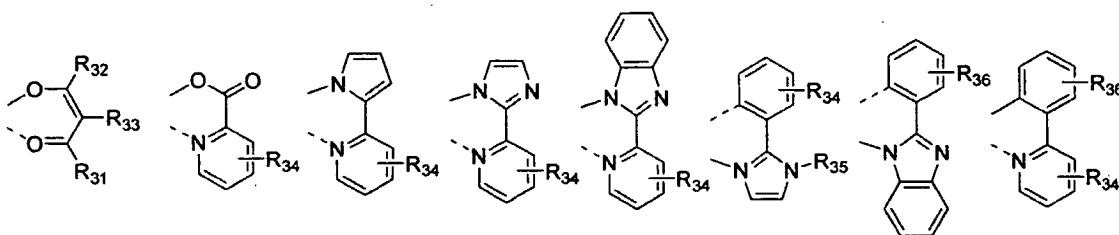
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R<sub>21</sub> represents (C1-C60)alkyl, halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, phenyl, di(C1-C60)alkylamino, di(C6-C60)arylamino, naphthyl, 9,9-di(C1-C60)alkylfluorenyl or 9,9-di(C6-C60)arylfluorenyl; and m is an integer from 1 to 5.

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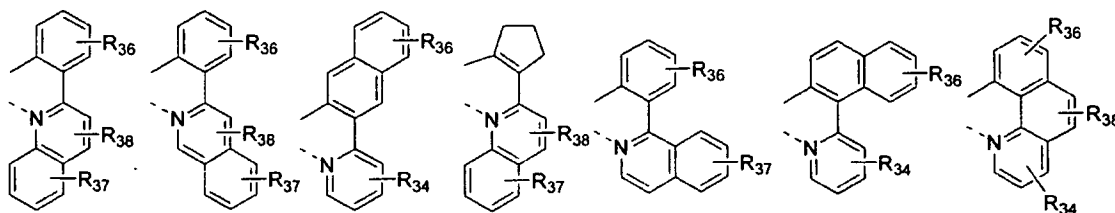
3. The organic electroluminescent compound according to claim 1, wherein the ligand (L) has a structure represented by one of the following chemical formulas:

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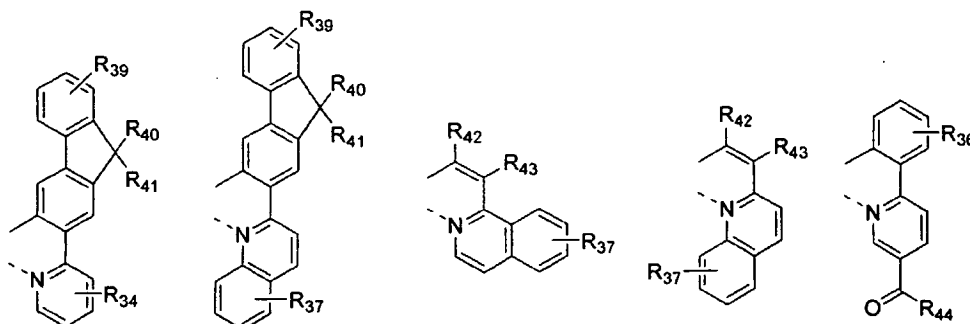
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wherein,  $R_{31}$  and  $R_{32}$  independently represent hydrogen, (C1-C60)alkyl with or without halogen substituent(s), phenyl with or without (C1-C60)alkyl substituent(s), or halogen;

$R_{33}$  through  $R_{39}$  independently represent hydrogen, (C1-C60)alkyl, phenyl with or without (C1-C60)alkyl substituent(s), tri(C1-C60)alkylsilyl or halogen;

20

$R_{40}$  through  $R_{43}$  independently represent hydrogen, (C1-C60)alkyl, phenyl with or without (C1-C60)alkyl substituent(s); and

$R_{44}$  represents (C1-C60)alkyl, phenyl with or without (C1-C60)alkyl substituent(s), or halogen.

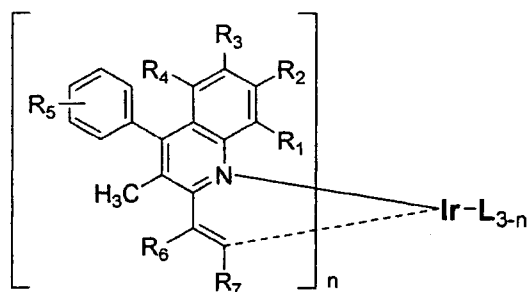
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4. An organic electroluminescent device which is comprised of a first electrode; a second electrode; and at least one organic layer(s) interposed between the first electrode and the second electrode; wherein the organic layer comprises an electroluminescent region comprising one or more compound(s) represented by an organic electroluminescent compound represented by Chemical Formula (1):

30

Chemical Formula 1

35



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wherein, L is an organic ligand;

$R_1$  through  $R_4$  independently represent hydrogen, (C1-C60)alkyl, (C1-C60)alkoxy, (C3-C60)cycloalkyl, halogen, tri(C1-C60)alkylsilyl or tri(C6-C60)arylsilyl;

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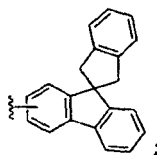
$R_5$  represents hydrogen, (C1-C60)alkyl, halogen, or (C6-C60)aryl;

$R_6$  and  $R_7$  independently represent hydrogen, (C1-C60)alkyl with or without halogen substituent(s), (C6-C60)aryl, halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino or di(C6-C60)arylamino, or  $R_6$  and  $R_7$  may be linked via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

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the alkyl or aryl of  $R_6$  and  $R_7$ , or the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed therefrom by linkage via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring may be further substituted by one or more substituent(s) selected from (C1-C60)alkyl with or without halogen substituent(s), halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino, di(C6-C60)arylamino, phenyl, naphthyl, anthryl, fluorenyl, spirobifluorenyl and

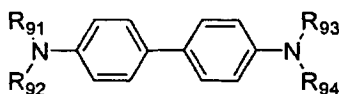
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and

n is an integer from 1 to 3, and one or more host(s) selected from 1,3,5-tricarbazolylbenzene, polyvinylcarbazole, m-biscarbazolylphenyl, 4,4'4''-tri(N-carbazolyl)triphenylamine, 1,3,5-tri(2-carbazolylphenyl)benzene, 1,3,5-tris(2-carbazolyl-5-methoxyphenyl)benzene, bis(4-carbazolylphenyl)silane and one or more host(s) selected from the compounds represented by one of Chemical Formulas (8) to (11):

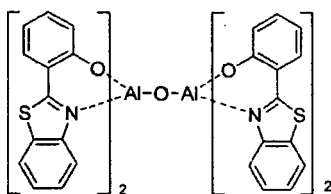
Chemical Formula 8



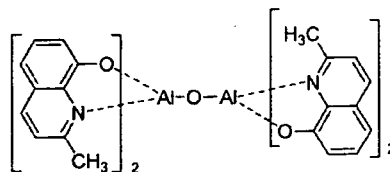
In Chemical Formula (8),  $R_{91}$  through  $R_{94}$  independently represent hydrogen, halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcabonyl, carboxyl, nitro or hydroxyl, or each of  $R_{91}$  through  $R_{94}$  may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

the alkyl, alkenyl, alkynyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, arylsilyl, alkylsilyl, alkylamino, or arylamino of  $R_{91}$  through  $R_{94}$ , or the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed therefrom by linkage to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring may be further substituted by one or more substituent(s) selected from halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcabonyl, carboxyl, nitro and hydroxyl.

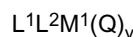
Chemical Formula 9



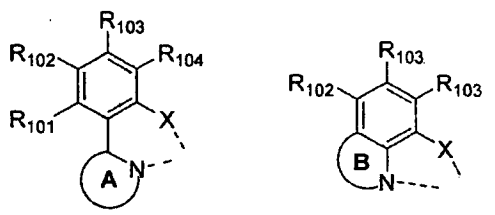
Chemical Formula 10



Chemical Formula 11



In Chemical Formula (11), the ligands,  $L^1$  and  $L^2$  are independently selected from the following structures:



$M^1$  is a bivalent or trivalent metal;

$y$  is 0 when  $M^1$  is a bivalent metal, while  $y$  is 1 when  $M^1$  is a trivalent metal;

$Q$  represents (C6-C60)aryloxy or tri(C6-C60)arylsilyl, and the aryloxy and triarylsilyl of  $Q$  may be further substituted by (C1-C60)alkyl or (C6-C60)aryl;

$X$  represents O, S or Se;

ring A represents oxazole, thiazole, imidazole, oxadiazole, thiadiazole, benzoxazole, benzothiazole, benzimidazole, pyridine or quinoline;

ring B represents pyridine or quinoline, and ring B may be further substituted by (C1-C60)alkyl, or phenyl or naphthyl with or without (C1-C60)alkyl substituent(s);

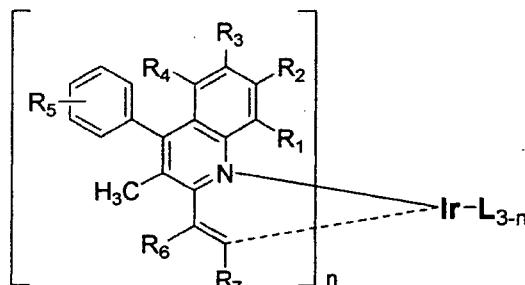
$R_{101}$  through  $R_{104}$  independently represent hydrogen, halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, carboxyl, nitro or hydroxyl, or each of  $R_{101}$  through  $R_{104}$  may be linked to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring; and the pyridine or quinoline may form a chemical bond with  $R_{101}$  to form a fused ring;

the alkyl, alkenyl, alkynyl, cycloalkyl, heterocycloalkyl, aryl, heteroaryl, arylsilyl, alkylsilyl, alkylamino, or arylamino of ring A and  $R_{101}$  through  $R_{104}$ , or the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed therefrom by linkage to an adjacent substituent via (C3-C60)alkylene or (C3-C60)alkenylene with or without a fused ring may be further substituted by one or more substituent(s) selected from halogen, (C1-C60)alkyl, (C6-C60)aryl, (C4-C60)heteroaryl, a 5- or 6-membered heterocycloalkyl containing one or more heteroatom(s) selected from N, O and S, (C3-C60)cycloalkyl, tri(C1-C60)alkylsilyl, di(C1-C60)alkyl(C6-C60)arylsilyl, tri(C6-C60)arylsilyl, adamantyl, (C7-C60)bicycloalkyl, (C2-C60)alkenyl, (C2-C60)alkynyl, cyano, (C1-C60)alkylamino, (C6-C60)arylamino, (C6-C60)ar(C1-C60)alkyl, (C1-C60)alkyloxy, (C1-C60)alkylthio, (C6-C60)aryloxy, (C6-C60)arylthio, (C1-C60)alkoxycarbonyl, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, carboxyl, nitro and hydroxyl.

5. The organic electroluminescent device according to claim 4, wherein the organic layer comprises one or more compound(s) selected from a group consisting of arylamine compounds and styrylamine compounds, or one or more metal(s) selected from a group consisting of organic metals of Group 1, Group 2, 4<sup>th</sup> period and 5<sup>th</sup> period transition metals, lanthanide metals and d-transition elements.

6. The organic electroluminescent device according to claim 4, which is an organic display comprising an organic electroluminescent compound represented by an organic electroluminescent compound represented by Chemical Formula (1) :

Chemical Formula 1



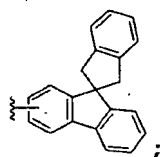
wherein, L is an organic ligand;

R<sub>1</sub> through R<sub>4</sub> independently represent hydrogen, (C1-C60)alkyl, (C1-C60)alkoxy, (C3-C60)cycloalkyl, halogen, tri(C1-C60)alkylsilyl or tri(C6-C60)arylsilyl;

R<sub>5</sub> represents hydrogen, (C1-C60)alkyl, halogen, or (C6-C60)aryl;

R<sub>6</sub> and R<sub>7</sub> independently represent hydrogen, (C1-C60)alkyl with or without halogen substituent(s), (C6-C60)aryl, halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino or di(C6-C60)arylamino, or R<sub>6</sub> and R<sub>7</sub> may be linked via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

the alkyl or aryl of R<sub>6</sub> and R<sub>7</sub>, or the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed therefrom by linkage via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring may be further substituted by one or more substituent(s) selected from (C1-C60)alkyl with or without halogen substituent(s), halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcarbonyl, di(C1-C60)alkylamino, di(C6-C60)arylamino, phenyl, naphthyl, anthryl, fluorenyl, spirobifluorenyl and

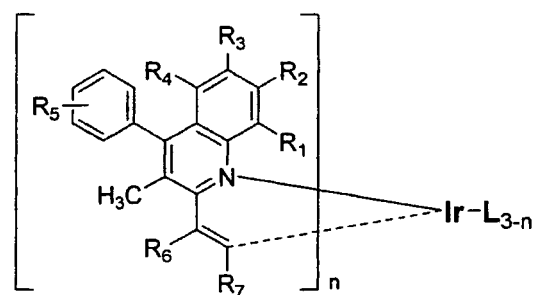


and

n is an integer from 1 to 3 and compounds having the electroluminescent peak with wavelength of blue and green at the same time.

7. The organic electroluminescent device according to claim 4, wherein the organic layer comprises an electroluminescent layer and a charge generating layer.
8. The organic electroluminescent device according to claim 4, wherein a mixed region of reductive dopant and organic substance, or a mixed region of oxidative dopant and organic substance is placed on the inner surface of one or both electrode(s) among the pair of electrodes.
9. An organic solar cell which comprises an organic electroluminescent compound represented by an organic electroluminescent compound represented by Chemical Formula (1) :

Chemical Formula 1



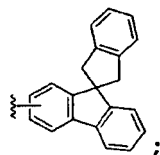
15 wherein, L is an organic ligand;

$R_1$  through  $R_4$  independently represent hydrogen, (C1-C60)alkyl, (C1-C60)alkoxy, (C3-C60)cycloalkyl, halogen, tri(C1-C60)alkylsilyl or tri(C6-C60)arylsilyl;

$R_5$  represents hydrogen, (C1-C60)alkyl, halogen, or (C6-C60)aryl;

20  $R_6$  and  $R_7$  independently represent hydrogen, (C1-C60)alkyl with or without halogen substituent(s), (C6-C60)aryl, halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcabonyl, di(C1-C60)alkylamino or di(C6-C60)arylamino, or  $R_6$  and  $R_7$  may may be linked via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring to form an alicyclic ring, or a monocyclic or polycyclic aromatic ring;

25 the alkyl or aryl of  $R_6$  and  $R_7$ , or the alicyclic ring, or the monocyclic or polycyclic aromatic ring formed therefrom by linkage via (C3-C12)alkylene or (C3-C12)alkenylene with or without a fused ring may be further substituted by one or more substituent(s) selected from (C1-C60)alkyl with or without halogen substituent(s), halogen, cyano, tri(C1-C60)alkylsilyl, tri(C6-C60)arylsilyl, (C1-C60)alkoxy, (C1-C60)alkylcarbonyl, (C6-C60)arylcabonyl, di(C1-C60)alkylamino, di(C6-C60)arylamino, phenyl, naphthyl, anthryl, fluorenyl, spirobifluorenyl and



and

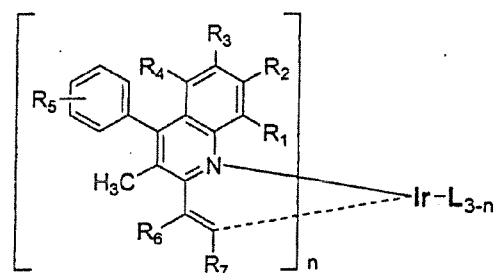
$n$  is an integer from 1 to 3.

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### Patentansprüche

- 45 1. Organische elektrolumineszente Verbindung, dargestellt durch die chemische Formel (1):

chemische Formel 1



wobei

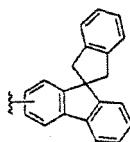
L ist ein organischer Ligand;

R<sub>1</sub> bis R<sub>4</sub> stellen unabhängig voneinander Wasserstoff, (C1-C60)Alkyl, (C1-C60)Alkoxy, (C3-C60)Cycloalkyl, Halogen, Tri(C1-C60)alkylsilyl oder Tri(C6-C60)arylsilyl dar;

R<sub>5</sub> stellt Wasserstoff, (C1-C60)Alkyl, Halogen oder (C6-C60)Aryl dar;

R<sub>6</sub> und R<sub>7</sub> stellen unabhängig voneinander Wasserstoff, (C1-C60)Alkyl mit oder ohne Halogensubstituent(en), (C6-C60)Aryl, Halogen, Cyano, Tri(C1-C60)alkylsilyl, Tri(C6-C60)arylsilyl, (C1-C60)Alkoxy, (C1-C60)Alkylcarbonyl, (C6-C60)Arylcarbonyl, Di(C1-C60)alkylamino oder Di(C6-C60)arylamino dar, oder R<sub>6</sub> und R<sub>7</sub> können über (C3-C12)Alkylen oder (C3-C12)Alkenylen mit oder ohne einen kondensierten Ring verbunden sein, um einen alicyclischen Ring, oder einen monocyclischen oder polycyclischen aromatischen Ring zu bilden;

das Alkyl oder Aryl von R<sub>6</sub> und R<sub>7</sub>, oder der davon über Bindung durch (C3-C12)Alkylen oder (C3-C12)Alkenylen mit oder ohne einen kondensierten Ring gebildete alicyclische Ring, oder der monocyclische oder polycyclische aromatische Ring kann weiter durch einen oder mehrere Substituenten, ausgewählt aus (C1-C60)Alkyl mit oder ohne Halogensubstituent(en), Halogen, Cyano, Tri(C1-C60)alkylsilyl, Tri(C6-C60)arylsilyl, (C1-C60)Alkoxy, (C1-C60)-Alkylcarbonyl, (C6-C60)Arylcarbonyl, Di(C1-C60)alkylamino, Di(C6-C60)arylamino, Phenyl, Naphthyl, Anthryl, Fluorenyl, Spirobifluorenyl und

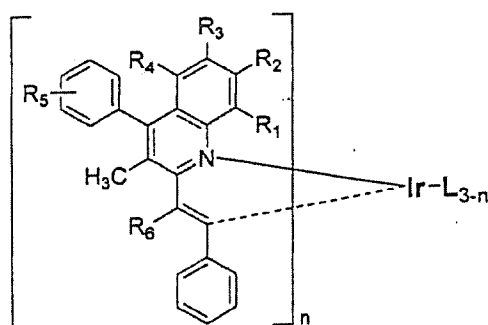


substituiert sein; und

n ist eine ganze Zahl von 1 bis 3.

2. Organische elektrolumineszente Verbindung nach Anspruch 1, die aus den Verbindungen, dargestellt durch eine der chemischen Formeln (2) bis (7), ausgewählt ist:

chemische Formel 2

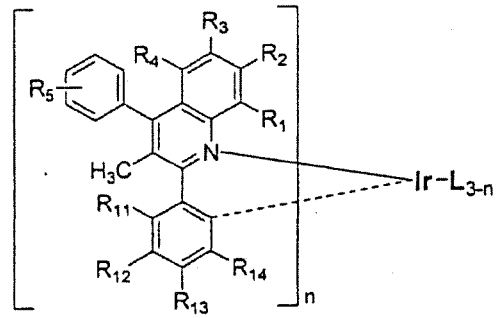


chemische Formel 3

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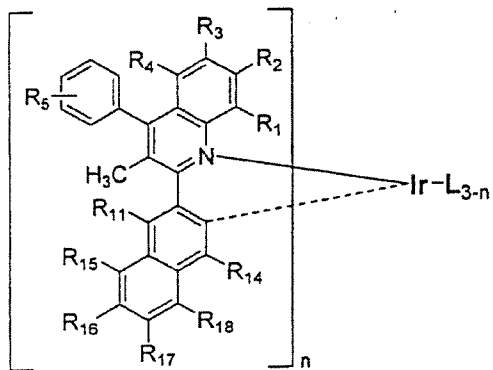


chemische Formel 4

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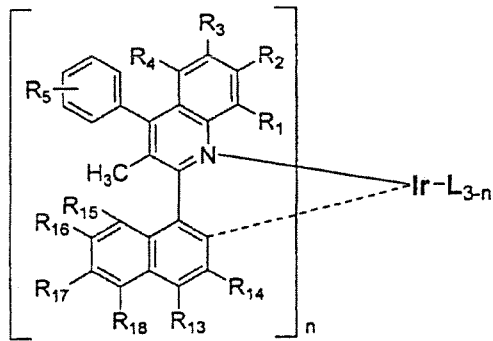


chemische Formel 5

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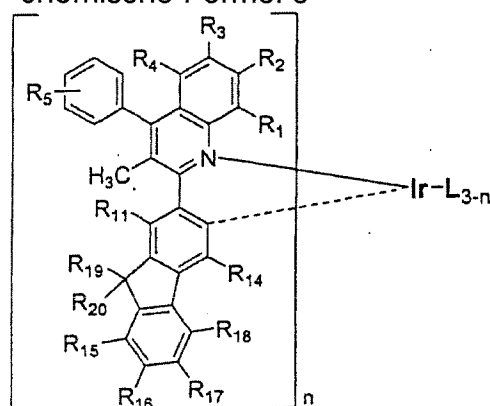
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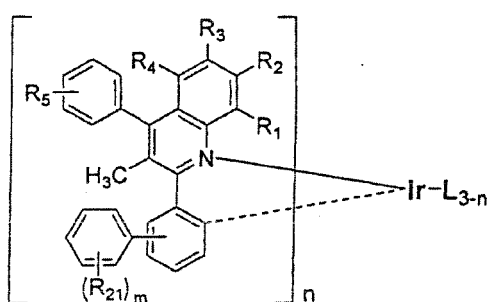
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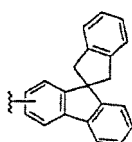
chemische Formel 6



chemische Formel 7



wobei L, R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub>, R<sub>5</sub>, R<sub>6</sub> und n wie in der chemischen Formel (1) von Anspruch 1 definiert sind; R<sub>11</sub> bis R<sub>18</sub> stellen unabhängig voneinander Wasserstoff, (C1-C60)Alkyl, Halogen, Cyano, Tri(C1-C60)alkylsilyl, Tri(C6-C60)arylsilyl, (C1-C60)Alkoxy, (C1-C60)Alkylcarbonyl, (C6-C60)Arylcarbonyl, Di(C1-C60)alkylamino, Di(C6-C60)arylamino, Phenyl, Naphthyl, Anthryl, Fluorenyl, Spirobifluorenyl oder



dar;

das Alkyl, Phenyl, Naphthyl, Anthryl oder Fluorenyl von R<sub>11</sub> bis R<sub>18</sub> kann weiter durch einen oder mehrere Substituenten, ausgewählt aus (C1-C60)Alkyl mit oder ohne Halogensubstituent(en), (C1-C60)Alkoxy, Halogen, Tri(C1-C60)alkylsilyl, Tri(C6-C60)arylsilyl, (C1-C60)Alkylcarbonyl, (C6-C60)Arylcarbonyl, Di(C1-C60)alkylamino, Di(C6-C60)arylamino und (C6-C60)Aryl substituiert sein;

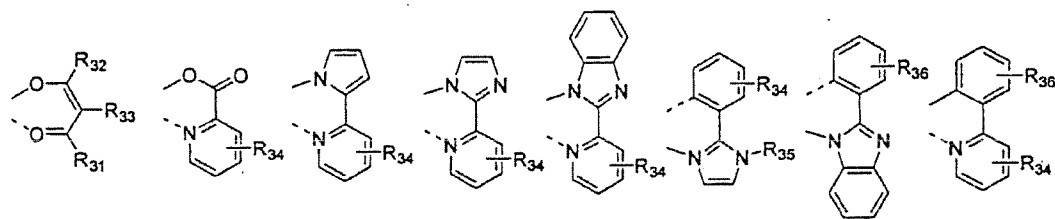
R<sub>19</sub> und R<sub>20</sub> stellen unabhängig voneinander Wasserstoff, (C1-C60)Alkyl oder (C6-C60)Aryl dar, oder R<sub>19</sub> und R<sub>20</sub> können über (C3-C12)Alkylen oder (C3-C12)Alkenylen mit oder ohne einen kondensierten Ring verbunden sein, um einen alicyclischen Ring, einen monocyclischen oder polycyclischen aromatischen Ring zu bilden;

R<sub>21</sub> stellt (C1-C60)Alkyl, Halogen, Cyano, Tri(C1-C60)alkylsilyl, Tri(C6-C60)arylsilyl, (C1-C60)Alkoxy, (C1-C60)Alkylcarbonyl, (C6-C60)Arylcarbonyl, Phenyl, Di(C1-C60)alkylamino, Di(C6-C60)arylamino, Naphthyl, 9,9-Di(C1-C60)alkylfluorenyl oder 9,9-Di(C6-C60)arylfluorenyl dar; und

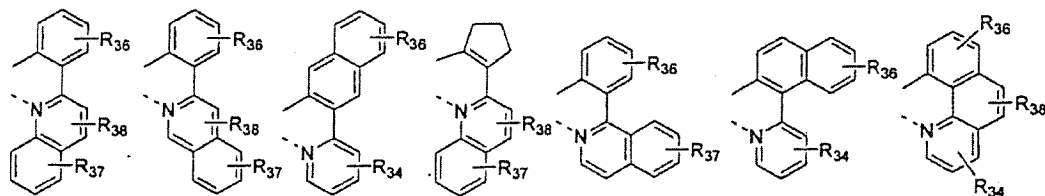
m ist eine ganze Zahl von 1 bis 5.

3. Organische elektrolumineszente Verbindung nach Anspruch 1, wobei der Ligand (L) eine durch eine der folgenden chemischen Formeln dargestellte Struktur aufweist:

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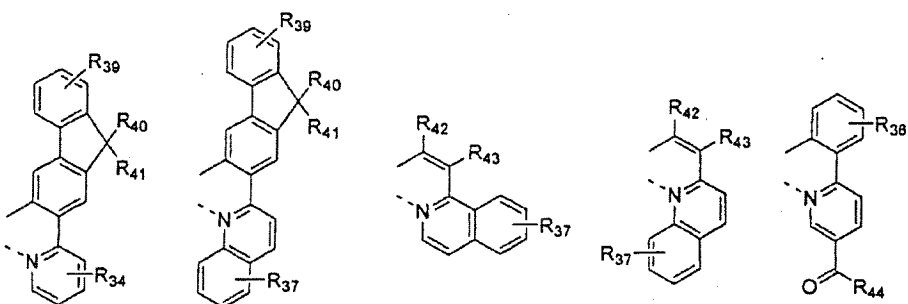
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wobei  $R_{31}$  und  $R_{32}$  unabhängig voneinander Wasserstoff, (C1-C60)Alkyl mit oder ohne Halogensubstituent(en), Phenyl mit oder ohne (C1-C60)Alkyl-substituent(en) oder Halogen darstellen;

35

$R_{33}$  bis  $R_{39}$  stellen unabhängig voneinander Wasserstoff, (C1-C60)Alkyl, Phenyl mit oder ohne (C1-C60)Alkylsubstituent(en), Tri(C1-C60)alkylsilyl oder Halogen dar;

$R_{40}$  bis  $R_{43}$  stellen unabhängig voneinander Wasserstoff, (C1-C60)Alkyl, Phenyl mit oder ohne (C1-C60)Alkylsubstituent(en) dar; und

$R_{44}$  stellt (C1-C60)Alkyl, Phenyl mit oder ohne (C1-C60)Alkylsubstituent(en) oder Halogen dar.

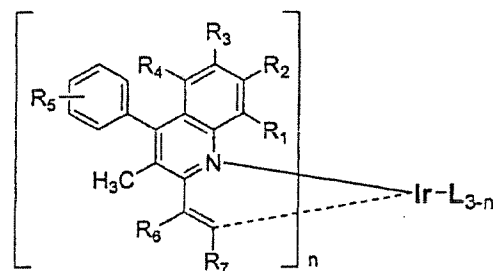
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4. Organische elektrolumineszente Vorrichtung, die eine erste Elektrode, eine zweite Elektrode und mindestens eine organische Schicht, welche zwischen der ersten Elektrode und der zweiten Elektrode angeordnet ist, umfasst, wobei die organische Schicht einen elektrolumineszenten Bereich, umfassend eine oder mehrere Verbindung(en), dargestellt durch eine organische elektrolumineszente Verbindung, die durch die chemische Formel (1) dargestellt ist:

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chemische Formel 1

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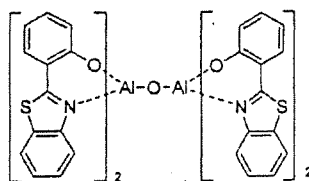


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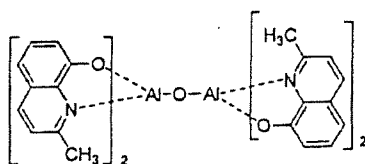


C60)Alkylcarbonyl, (C6-C60)Arylcarbonyl, Carboxyl, Nitro und Hydroxyl substituiert sein;

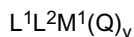
chemische Formel 9



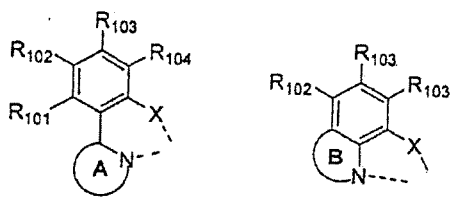
chemische Formel 10



chemische Formel 11



wobei in der chemischen Formel (11) die Liganden,  $L^1$  und  $L^2$ , unabhängig voneinander aus den folgenden Formeln ausgewählt sind:



$M^1$  ist ein zweiwertiges oder dreiwertiges Metall;

$y$  ist 0, falls  $M^1$  ein zweiwertiges Metall ist, wohingegen  $y$  1 ist, falls  $M^1$  ein dreiwertiges Metall ist;

$Q$  stellt (C6-C60)Aryloxy oder Tri(C6-C60)arylsilyl dar, und das Aryloxy und Triarylsilyl von  $Q$  kann weiter durch (C1-C60)Alkyl oder (C6-C60)Aryl substituiert sein;

$X$  stellt O, S oder Se dar;

Ring A stellt Oxazol, Thiazol, Imidazol, Oxadiazol, Thiadiazol, Benzoxazol, Benzothiazol, Benzimidazol, Pyridin oder Chinolin dar;

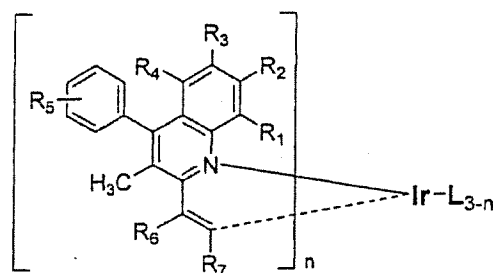
Ring B stellt Pyridin oder Chinolin dar, und Ring B kann weiter durch (C1-C60)Alkyl oder Phenyl oder Naphthyl mit oder ohne (C1-C60)Alkyl-substituent(en) substituiert sein;

$R_{101}$  bis  $R_{104}$  stellen unabhängig voneinander Wasserstoff, Halogen, (C1-C60)Alkyl, (C6-C60)Aryl, (C4-C60)Heteroaryl, ein 5- oder 6-gliedriges Heterocycloalkyl, das ein oder mehrere Heteroatom(e), ausgewählt aus N, O und S, enthält, (C3-C60)Cycloalkyl, Tri(C1-C60)alkylsilyl, Di(C1-C60)alkyl(C6-C60)arylsilyl, Tri(C6-C60)arylsilyl, Adamantyl, (C7-C60)Bicycloalkyl, (C2-C60)Alkenyl, (C2-C60)Alkyl, Cyano, (C1-C60)Alkylamino, (C6-C60)Arylamino, (C6-C60)Ar(C1-C60)alkyl, (C1-C60)Alkyl, (C1-C60)Alkylthio, (C6-C60)Aryloxy, (C6-C60)Arylthio, (C1-C60)Alkoxy, (C1-C60)Alkylcarbonyl, (C6-C60)Arylcarbonyl, Carboxyl, Nitro oder Hydroxyl dar, oder jedes von  $R_{101}$  bis  $R_{104}$  kann zu einem benachbarten Substituenten über (C3-C60)Alkylen oder (C3-C60)Alkenylen mit oder ohne einen kondensierten Ring verbunden sein, um einen alicyclischen Ring, oder einen monocyclischen oder polycyclischen aromatischen Ring zu bilden; und das Pyridin oder Chinolin kann

mit  $R_{101}$  eine chemische Bindung bilden, um einen kondensierten Ring zu bilden;  
 das Alkyl, Alkenyl, Alkynyl, Cycloalkyl, Heterocycloalkyl, Aryl, Heteroaryl, Arylsilyl, Alkylsilyl, Alkylamino oder  
 Arylamino von Ring A und  $R_{101}$  bis  $R_{104}$ , oder der davon über Bindung zu einem benachbarten Substituenten  
 durch (C3-C60)Alkylen oder (C3-C60)Alkenylen mit oder ohne einen kondensierten Ring gebildete alicyclische  
 Ring, oder der monocyclische oder polycyclische aromatische Ring kann weiter durch einen oder mehrere  
 Substituenten, ausgewählt aus Halogen, (C1-C60)Alkyl, (C6-C60)Aryl, (C4-C60)Heteroaryl, einem 5- oder 6-  
 gliedrigen Heterocycloalkyl, das ein oder mehrere Heteroatom(e), ausgewählt aus N, O und S, enthält, (C3-  
 C60)Cycloalkyl, Tri(C1-C60)alkylsilyl, Di(C1-C60)alkyl(C6-C60)arylsilyl, Tri(C6-C60)arylsilyl, Adamantyl, (C7-  
 C60)Bicycloalkyl, (C2-C60)Alkenyl, (C2-C60)Alkynyl, Cyano, (C1-C60)Alkylamino, (C6-C60)Arylamino, (C6-  
 C60)Ar(C1-C60)alkyl, (C1-C60)Alkyloxy, (C1-C60)-Alkylthio, (C6-C60)Aryloxy, (C6-C60)Arylthio, (C1-C60)Alk-  
 oxycarbonyl, (C1-C60)Alkylcarbonyl, (C6-C60)Arylcarbonyl, Carboxyl, Nitro und Hydroxyl, substituiert sein.

5. Organische elektrolumineszente Vorrichtung nach Anspruch 4, wobei die organische Schicht eine oder mehrere  
 Verbindung(en), ausgewählt aus der Gruppe, bestehend aus Arylaminverbindungen und Styrylaminverbindungen,  
 oder ein oder mehrere Metall(e), ausgewählt aus der Gruppe, bestehend aus organischen Metallen der Gruppe 1,  
 Gruppe 2, Übergangsmetallen der 4. Periode und der 5. Periode, Lanthanidmetallen und d-Übergangselementen,  
 umfasst.
6. Organische elektrolumineszente Vorrichtung nach Anspruch 4, die ein organisches Display ist, umfassend eine  
 organische elektrolumineszente Verbindung, dargestellt durch eine organische elektrolumineszente Verbindung,  
 die durch die chemische Formel (1) dargestellt ist:

chemische Formel 1



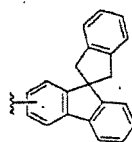
wobei

L ist ein organischer Ligand;

$R_1$  bis  $R_4$  stellen unabhängig voneinander Wasserstoff, (C1-C60)Alkyl, (C1-C60)Alkoxy, (C3-C60)Cycloalkyl,  
 Halogen, Tri(C1-C60)alkylsilyl oder Tri(C6-C60)arylsilyl dar;

$R_5$  stellt Wasserstoff, (C1-C60)Alkyl, Halogen oder (C6-C60)Aryl dar;

$R_6$  und  $R_7$  stellen unabhängig voneinander Wasserstoff, (C1-C60)Alkyl mit oder ohne Halogensubstituent(en),  
 (C6-C60)Aryl, Halogen, Cyano, Tri(C1-C60)alkylsilyl, Tri(C6-C60)arylsilyl, (C1-C60)Alkoxy, (C1-C60)Alkylcar-  
 bonyl, (C6-C60)Arylcarbonyl, Di(C1-C60)alkylamino oder Di(C6-C60)arylamino dar, oder  $R_6$  und  $R_7$  können  
 über (C3-C12)Alkylen oder (C3-C12)Alkenylen mit oder ohne einen kondensierten Ring verbunden sein, um  
 einen alicyclischen Ring, oder einen monocyclischen oder polycyclischen aromatischen Ring zu bilden;  
 das Alkyl oder Aryl von  $R_6$  und  $R_7$ , oder der davon über Bindung durch (C3-C12)Alkylen oder (C3-C12)Alkenylen  
 mit oder ohne einen kondensierten Ring gebildete alicyclische Ring, oder der monocyclische oder polycyclische  
 aromatische Ring kann weiter durch einen oder mehrere Substituenten, ausgewählt aus (C1-C60)Alkyl mit oder  
 ohne Halogensubstituent(en), Halogen, Cyano, Tri(C1-C60)alkylsilyl, Tri(C6-C60)arylsilyl, (C1-C60)Alkoxy,  
 (C1-C60)Alkylcarbonyl, (C6-C60)Arylcarbonyl, Di(C1-C60)alkylamino, Di(C6-C60)arylamino, Phenyl, Naphthyl,  
 Anthryl, Fluorenyl, Spirobifluorenyl und



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substituiert sein; und n ist eine ganze Zahl von 1 bis 3, und Verbindungen, welche den Elektrolumineszenzpeak gleichzeitig bei Wellenlängen von Blau und Grün aufweisen.

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7. Organische elektrolumineszente Vorrichtung nach Anspruch 4, wobei die organische Schicht eine Elektrolumineszenzschicht und eine Ladungserzeugungsschicht umfasst.

8. Organische elektrolumineszente Vorrichtung nach Anspruch 4, wobei ein gemischter Bereich aus reduzierendem Dotierstoff und organischer Substanz, oder ein gemischter Bereich aus oxidierendem Dotierstoff und organischer Substanz auf der Innenfläche von einer oder beiden Elektrode(n) zwischen dem Elektrodenpaar angeordnet ist.

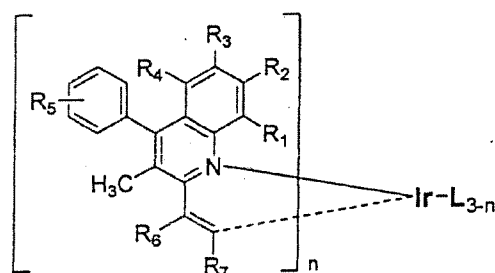
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9. Organische Solarzelle, die eine organische elektrolumineszente Verbindung umfasst, dargestellt durch eine organische elektrolumineszente Verbindung, die durch die chemische Formel (1) dargestellt ist:

## chemische Formel 1

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wobei

L ist ein organischer Ligand;

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R<sub>1</sub> bis R<sub>4</sub> stellen unabhängig voneinander Wasserstoff, (C1-C60)Alkyl, (C1-C60)Alkoxy, (C3-C60)Cycloalkyl, Halogen, Tri(C1-C60)alkylsilyl oder Tri(C6-C60)arylsilyl dar;

R<sub>5</sub> stellt Wasserstoff, (C1-C60)Alkyl, Halogen oder (C6-C60)Aryl dar;

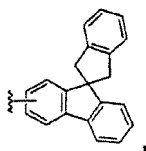
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R<sub>6</sub> und R<sub>7</sub> stellen unabhängig voneinander Wasserstoff, (C1-C60)Alkyl mit oder ohne Halogensubstituent(en), (C6-C60)Aryl, Halogen, Cyano, Tri(C1-C60)alkylsilyl, Tri(C6-C60)arylsilyl, (C1-C60)Alkoxy, (C1-C60)Alkylcarbonyl, (C6-C60)Arylcarbonyl, Di(C1-C60)alkylamino oder Di(C6-C60)arylamino dar, oder R<sub>6</sub> und R<sub>7</sub> können über (C3-C12)Alkylen oder (C3-C12)Alkenylen mit oder ohne einen kondensierten Ring verbunden sein, um einen alicyclischen Ring, oder einen monocyclischen oder polycyclischen aromatischen Ring zu bilden;

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das Alkyl oder Aryl von R<sub>6</sub> und R<sub>7</sub>, oder der davon über Bindung durch (C3-C12)Alkylen oder (C3-C12)Alkenylen mit oder ohne einen kondensierten Ring gebildete alicyclische Ring, oder der monocyclische oder polycyclische aromatische Ring kann weiter durch einen oder mehrere Substituenten, ausgewählt aus (C1-C60)Alkyl mit oder ohne Halogensubstituent(en), Halogen, Cyano, Tri(C1-C60)alkylsilyl, Tri(C6-C60)arylsilyl, (C1-C60)Alkoxy, (C1-C60)Alkylcarbonyl, (C6-C60)Arylcarbonyl, Di(C1-C60)alkylamino, Di(C6-C60)-arylamino, Phenyl, Naphthyl, Anthryl, Fluorenyl, Spirofluorenyl und

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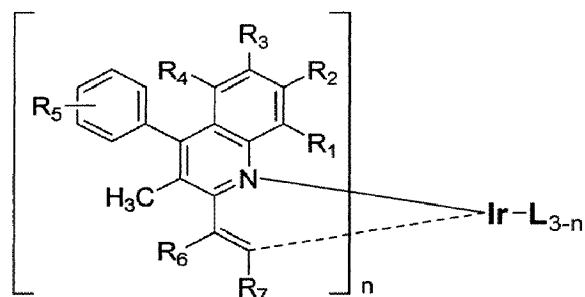


substitué(e) sein; und  
n ist eine ganze Zahl von 1 bis 3.

## Revendications

1. Composé électroluminescent organique représenté par la formule chimique (1) :

### formule chimique 1



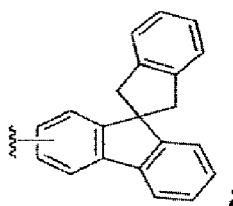
dans laquelle, L est un ligand organique ;

R<sub>1</sub> à R<sub>4</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, alcoxy en C<sub>1</sub>-C<sub>60</sub>, cycloalkyle en C<sub>3</sub>-C<sub>60</sub>, un halogène, un groupe tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle ou tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle ;

R<sub>5</sub> représente l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, un halogène ou un groupe aryle en C<sub>6</sub>-C<sub>60</sub> ;

R<sub>6</sub> et R<sub>7</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> avec ou sans substituant(s) d'halogène, un groupe aryle en C<sub>6</sub>-C<sub>60</sub>, un halogène, un groupe cyano, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, alcoxy en C<sub>1</sub>-C<sub>60</sub>, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino ou di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino, ou R<sub>6</sub> et R<sub>7</sub> peuvent être liés via un groupe alkylène en C<sub>3</sub>-C<sub>12</sub> ou alcénylène en C<sub>3</sub>-C<sub>12</sub> avec ou sans noyau condensé pour former un noyau alicyclique, ou un noyau aromatique monocyclique ou polycyclique ;

le groupe alkyle ou aryle de R<sub>6</sub> et R<sub>7</sub>, ou le noyau alicyclique, ou le noyau aromatique monocyclique ou polycyclique formé à partir de ceux-ci par liaison via un groupe alkylène en C<sub>3</sub>-C<sub>12</sub> ou alcénylène en C<sub>3</sub>-C<sub>12</sub> avec ou sans noyau condensé peuvent encore être substitués par un ou plusieurs substituant(s) choisi(s) parmi un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> avec ou sans substituant(s) d'halogène, un halogène, un groupe cyano, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, alcoxy en C<sub>1</sub>-C<sub>60</sub>, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino, di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino, phényle, naphtyle, anthryle, fluorényle, spirobi-fluorényle et

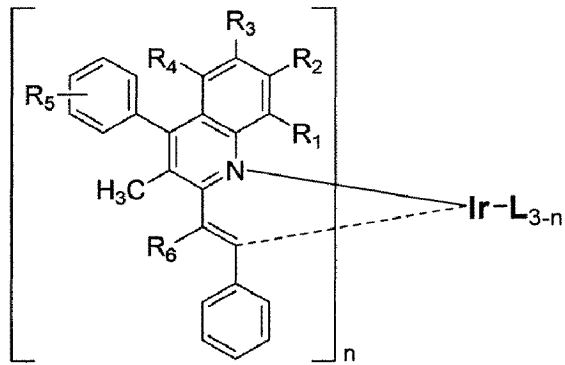


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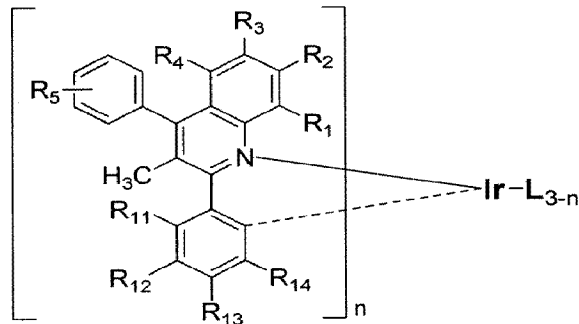
n est un nombre entier de 1 à 3.

2. Composé électroluminescent organique selon la revendication 1, lequel est choisi parmi les composés représentés par l'une des formules chimiques (2) à (7) :

formule chimique 2



formule chimique 3



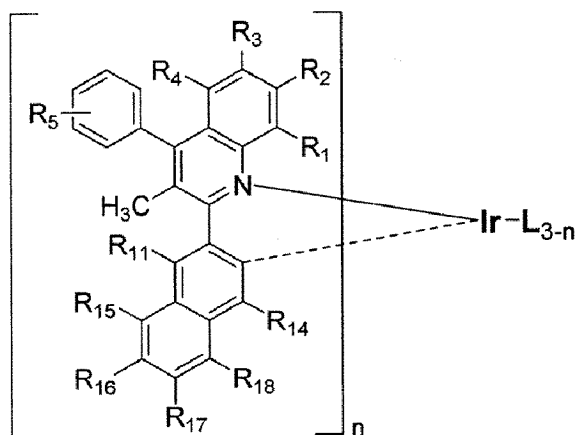
formule chimique 4

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formule chimique 5

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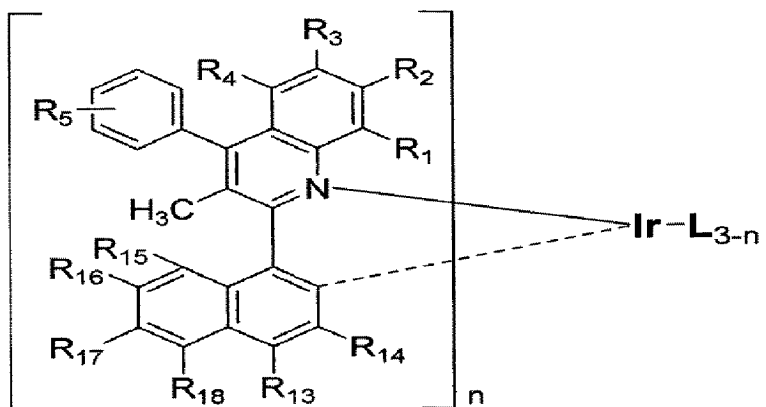
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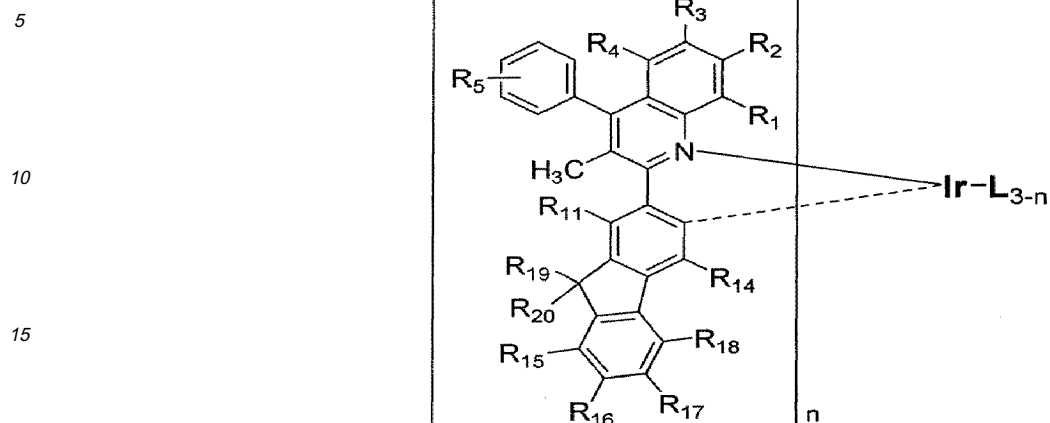
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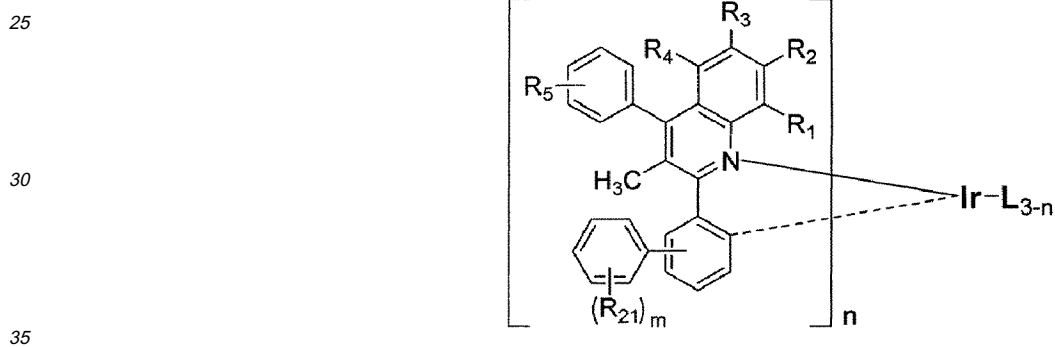


formule chimique 6



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formule chimique 7

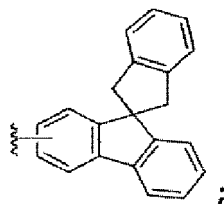


dans lesquelles L, R<sub>1</sub>, R<sub>2</sub>, R<sub>3</sub>, R<sub>4</sub>, R<sub>5</sub>, R<sub>6</sub> et n sont comme définis dans la formule chimique (1) de la revendication 1 ;

40 R<sub>11</sub> à R<sub>18</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, un halogène, un groupe cyano, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, alcoxy en C<sub>1</sub>-C<sub>60</sub>, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino, di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino, phényle, naphtyle, anthryle, fluorényle, spirobifluorényle ou

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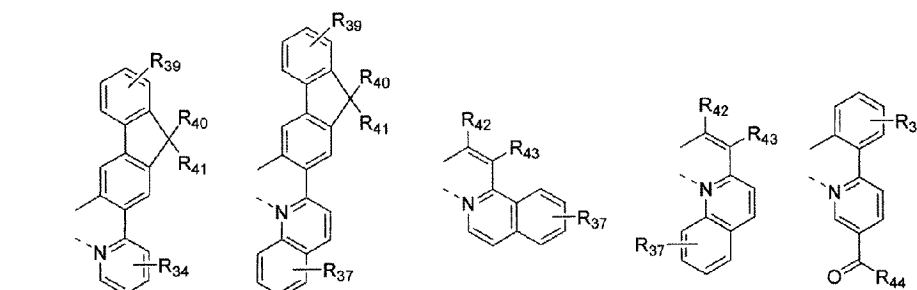
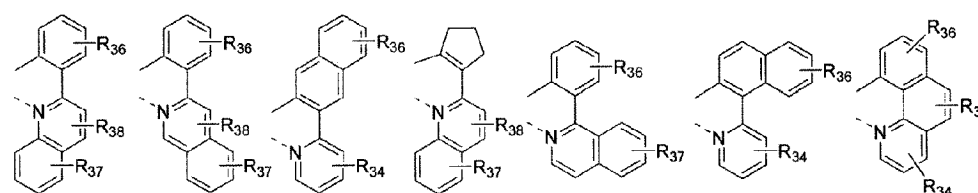
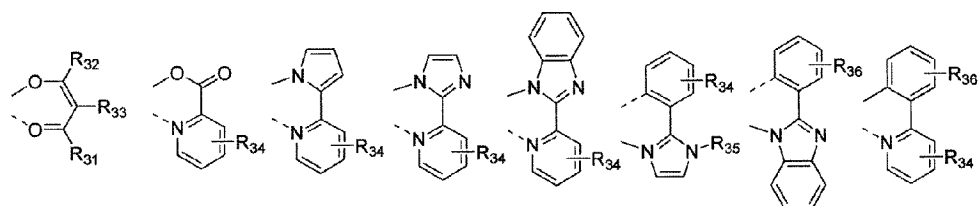
les groupes alkyle, phényle, naphtyle, anthryle ou fluorényle de R<sub>11</sub> à R<sub>18</sub> peuvent encore être substitués par un ou plusieurs substituant(s) choisi(s) parmi un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> avec ou sans substituant(s) d'halogène, un groupe alcoxy en C<sub>1</sub>-C<sub>60</sub>, un halogène, un groupe tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino, di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino et aryle en C<sub>6</sub>-C<sub>60</sub> ;

R<sub>19</sub> et R<sub>20</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> ou aryle en C<sub>6</sub>-C<sub>60</sub>, ou R<sub>19</sub> et R<sub>20</sub> peuvent être liés via un groupe alkylène en C<sub>3</sub>-C<sub>12</sub> ou alcénylène en C<sub>3</sub>-C<sub>12</sub> avec ou sans noyau condensé pour former un noyau alicyclique, un noyau aromatique monocyclique ou polycyclique ;

R<sub>21</sub> représente un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, un halogène, un groupe cyano, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, alcoxy en C<sub>1</sub>-C<sub>60</sub>, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, phényle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino, di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino, naphtyle, 9,9-di(alkyle en C<sub>1</sub>-C<sub>60</sub>)fluorényle ou 9,9-di(aryle en C<sub>6</sub>-C<sub>60</sub>)fluorényle ; et

m est un nombre entier de 1 à 5.

3. Composé électroluminescent organique selon la revendication 1, dans lequel le ligand (L) présente une structure représentée par l'une des formules chimiques suivantes :



dans lesquelles, R<sub>31</sub> et R<sub>32</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> avec ou sans substituant(s) d'halogène, phényle avec ou sans substituant(s) alkyle en C<sub>1</sub>-C<sub>60</sub>, ou un halogène ;

R<sub>33</sub> à R<sub>39</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, phényle avec ou sans substituant(s) alkyle en C<sub>1</sub>-C<sub>60</sub>, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle ou un halogène ;

R<sub>40</sub> à R<sub>43</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, phényle avec ou sans substituant(s) alkyle en C<sub>1</sub>-C<sub>60</sub> ; et

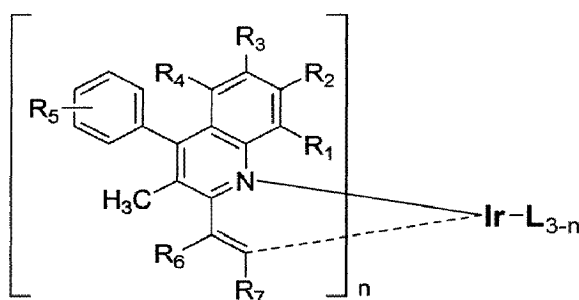
R<sub>44</sub> représente un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, phényle avec ou sans substituant(s) alkyle en C<sub>1</sub>-C<sub>60</sub> ou un halogène.

4. Dispositif électroluminescent organique qui est constitué d'une première électrode ; d'une seconde électrode ; et d'au moins une couche organique interposée entre la première électrode et la seconde électrode ; dans lequel la couche organique comprend une région électroluminescente comprenant un ou plusieurs composé(s) représenté(s) par un composé électroluminescent organique représenté par la formule chimique (I) :

## formule chimique 1

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dans laquelle, L est un ligand organique ;

R<sub>1</sub> à R<sub>4</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, alcoxy en C<sub>1</sub>-C<sub>60</sub>, cycloalkyle en C<sub>3</sub>-C<sub>60</sub>, un halogène, un groupe tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle ou tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle ;

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R<sub>5</sub> représente l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, un halogène ou un groupe aryle en C<sub>6</sub>-C<sub>60</sub> ;

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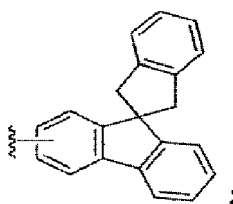
R<sub>6</sub> et R<sub>7</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> avec ou sans substituant(s) d'halogène, un groupe aryle en C<sub>6</sub>-C<sub>60</sub>, un halogène, un groupe cyano, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, alcoxy en C<sub>1</sub>-C<sub>60</sub>, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino ou di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino, ou R<sub>6</sub> et R<sub>7</sub> peuvent être liés via un groupe alkylène en C<sub>3</sub>-C<sub>12</sub> ou alcénylène en C<sub>3</sub>-C<sub>12</sub> avec ou sans noyau condensé pour former un noyau alicyclique, ou un noyau aromatique monocyclique ou polycyclique ;

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le groupe alkyle ou aryle de R<sub>6</sub> et R<sub>7</sub>, ou le noyau alicyclique, ou le noyau aromatique monocyclique ou polycyclique formé à partir de ceux-ci par liaison via un groupe alkylène en C<sub>3</sub>-C<sub>12</sub> ou alcénylène en C<sub>3</sub>-C<sub>12</sub> avec ou sans noyau condensé peuvent encore être substitués par un ou plusieurs substituant(s) choisi(s) parmi un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> avec ou sans substituant(s) d'halogène, un halogène, un groupe cyano, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, alcoxy en C<sub>1</sub>-C<sub>60</sub>, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino, di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino, phényle, naphtyle, anthryle, fluorényle, spirobi-fluorényle et

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et

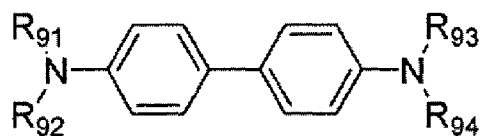
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n est un nombre entier de 1 à 3, et un ou plusieurs hôte(s) choisi(s) parmi le 1,3,5-tricarbazolylbenzène, le polyvinylcarbazole, le m-biscarbazolylphényle, la 4,4',4''-tri(N-carbazolyl)triphénylamine, le 1,3,5-tri(2-carbazolylphényl)benzène, le 1,3,5-tris(2-carbazolyl-5-méthoxyphényl)benzène, le bis(4-carbazolylphényl)silane et un ou plusieurs hôte(s) choisi(s) parmi les composés représentés par l'une des formules chimiques (8) à (11) :

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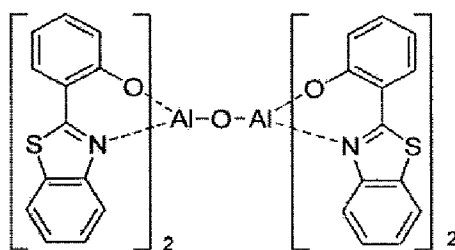
## formule chimique 8



dans la formule chimique (8),  $R_{91}$  à  $R_{94}$  représentent indépendamment l'hydrogène, un halogène, un groupe alkyle en  $C_1-C_{60}$ , aryle en  $C_6-C_{60}$ , hétéroaryle en  $C_4-C_{60}$ , hétérocycloalkyle à 5 ou 6 éléments contenant un ou plusieurs hétéroatome(s) choisi(s) parmi N, O et S, cycloalkyle en  $C_3-C_{60}$ , tri(alkyle en  $C_1-C_{60}$ )silyle, di(alkyle en  $C_1-C_{60}$ ) (aryle en  $C_6-C_{60}$ )silyle, tri(aryle en  $C_6-C_{60}$ )silyle, adamantyle, bicycloalkyle en  $C_7-C_{60}$ , alcényle en  $C_2-C_{60}$ , alcynyle en  $C_2-C_{60}$ , cyano, (alkyle en  $C_1-C_{60}$ )amino, (aryle en  $C_6-C_{60}$ )amino, (aryle en  $C_6-C_{60}$ ) (alkyle en  $C_1-C_{60}$ ), (alkyle en  $C_1-C_{60}$ )oxy, (alkyle en  $C_1-C_{60}$ )thio, (aryle en  $C_6-C_{60}$ )oxy, (aryle en  $C_6-C_{60}$ )thio, (alcoxy en  $C_1-C_{60}$ )carbonyle, (alkyle en  $C_1-C_{60}$ )carbonyle, (aryle en  $C_6-C_{60}$ )carbonyle, carboxyle, nitro ou hydroxyle ou chacun de  $R_{91}$  à  $R_{94}$  peut être lié à un substituant adjacent via un groupe alkylène en  $C_3-C_{60}$  ou alcénylène en  $C_3-C_{60}$  avec ou sans noyau condensé pour former un noyau alicyclique ou un noyau aromatique monocyclique ou polycyclique ;

le groupe alkyle, alcényle, alcynyle, cycloalkyle, hétérocycloalkyle, aryle, hétéroaryle, arylsilyle, alkylsilyle, alkylamino ou arylamino de  $R_{91}$  à  $R_{94}$ , ou le noyau alicyclique, ou le noyau aromatique monocyclique ou polycyclique formé à partir de ceux-ci par liaison à un substituant adjacent via un groupe alkylène en  $C_3-C_{60}$  ou alcénylène en  $C_3-C_{60}$  avec ou sans noyau condensé peuvent encore être substitués par un ou plusieurs substituant(s) choisi(s) parmi un halogène, un groupe alkyle en  $C_1-C_{60}$ , aryle en  $C_6-C_{60}$ , hétéroaryle en  $C_4-C_{60}$ , hétérocycloalkyle à 5 ou 6 éléments contenant un ou plusieurs hétéroatome(s) choisi(s) parmi N, O et S, cycloalkyle en  $C_3-C_{60}$ , tri(alkyle en  $C_1-C_{60}$ )silyle, di(alkyle en  $C_1-C_{60}$ ) (aryle en  $C_6-C_{60}$ )silyle, tri(aryle en  $C_6-C_{60}$ )silyle, adamantyle, bicycloalkyle en  $C_7-C_{10}$ , alcényle en  $C_2-C_{60}$ , alcynyle en  $C_2-C_{60}$ , cyano, (alkyle en  $C_1-C_{60}$ )amino, (aryle en  $C_6-C_{60}$ )amino, (aryle en  $C_6-C_{60}$ ) (alkyle en  $C_1-C_{60}$ ), (alkyle en  $C_1-C_{60}$ )oxy, (alkyle en  $C_1-C_{60}$ )thio, (aryle en  $C_6-C_{60}$ )oxy, (aryle en  $C_6-C_{60}$ )thio, (alcoxy en  $C_1-C_{60}$ )carbonyle, (alkyle en  $C_1-C_{60}$ )carbonyle, (aryle en  $C_6-C_{60}$ )carbonyle, carboxyle, nitro et hydroxyle.

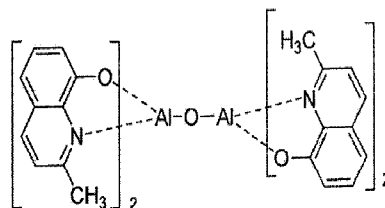
## Formule chimique 9



## Formule chimique 10

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Formule chimique 11

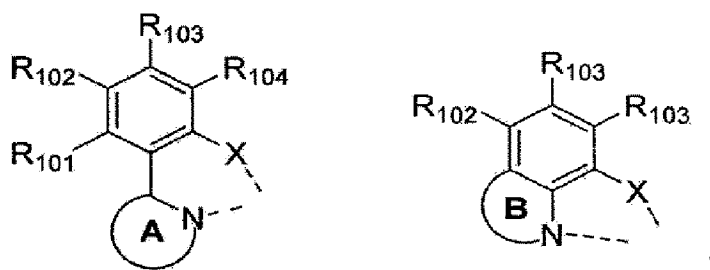
 $L^1L^2M^1(Q)_y$ 

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dans la formule chimique (11), les ligands,  $L^1$  et  $L^2$  sont indépendamment choisis parmi les structures suivantes :

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$M^1$  est un métal bivalent ou trivalent ;

$y$  est égal à 0 lorsque  $M^1$  est un métal bivalent alors que  $y$  est égal à 1 lorsque  $M^1$  est un métal trivalent ;

$Q$  représente un groupe (aryle en  $C_6-C_{60}$ )oxy ou tri(aryle en  $C_6-C_{60}$ )silyle, et les groupes aryloxy et triarylsilyle de  $Q$  peuvent encore être substitués par un groupe alkyle en  $C_1-C_{60}$  ou aryle en  $C_6-C_{60}$  ;

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$X$  représente O, S ou Se ;

le noyau A représente un noyau oxazole, thiazole, imidazole, oxadiazole, thiadiazole, benzoxazole, benzothiazole, benzimidazole, pyridine ou quinoléine ;

le noyau B représente un noyau pyridine ou quinoléine, et le noyau B peut encore être substitué par un groupe alkyle en  $C_1-C_{60}$ , ou phényle ou naphtyle avec ou sans substituant(s) alkyle en  $C_1-C_{60}$  ;

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$R_{101}$  à  $R_{104}$  représentent indépendamment l'hydrogène, un halogène, un groupe alkyle en  $C_1-C_{60}$ , aryle en  $C_6-C_{60}$ , hétéroaryle en  $C_4-C_{60}$ , hétérocycloalkyle à 5 ou 6 chaînons contenant un ou plusieurs hétéroatome(s) choisi(s) parmi N, O et S, cycloalkyle en  $C_3-C_{60}$ , tri(alkyle en  $C_1-C_{60}$ )Silyle, di(alkyle en  $C_1-C_{60}$ ) (aryle en  $C_6-C_{60}$ ) silyle, tri(aryle en  $C_6-C_{60}$ )silyle, adamantyle, bicycloalkyle en  $C_7-C_{60}$ , alcényle en  $C_2-C_{60}$ , alcynyle en  $C_2-C_{60}$ , cyano, (alkyle en  $C_1-C_{60}$ )amino, (aryle en  $C_6-C_{60}$ )amino, (aryle en  $C_6-C_{60}$ ) (alkyle en  $C_1-C_{60}$ ), (alkyle en  $C_1-C_{60}$ ) oxy, (alkyle en  $C_1-C_{60}$ )thio, (aryle en  $C_6-C_{60}$ )oxy, (aryle en  $C_6-C_{60}$ )thio, (alcoxy en  $C_1-C_{60}$ )carbonyle, (alkyle en  $C_1-C_{60}$ )carbonyle, (aryle en  $C_6-C_{60}$ )carbonyle, carboxyle, nitro ou hydroxyle, ou chacun de  $R_{101}$  à  $R_{104}$  peut être lié à un substituant adjacent via un groupe alkylène en  $C_3-C_{60}$  ou alcénylène en  $C_3-C_{60}$  avec ou sans noyau condensé pour former un noyau alicyclique ou un noyau aromatique monocyclique ou polycyclique ; et

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le noyau pyridine ou quinoléine peut former une liaison chimique avec  $R_{101}$  pour former un noyau condensé ;

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les groupes alkyle, alcényle, alcynyle, cycloalkyle, hétérocycloalkyle, aryle, hétéroaryle, arylsilyle, alkylsilyle, alkylamino ou arylamino du noyau A et  $R_{101}$  à  $R_{104}$ , ou le noyau alicyclique ou le noyau aromatique monocyclique ou polycyclique formé à partir de ceux-ci par liaison à un substituant adjacent via un groupe alkylène en  $C_3-C_{60}$  ou alcénylène en  $C_3-C_{60}$  avec ou sans noyau condensé peuvent encore être substitués par un ou plusieurs substituant(s) choisi(s) parmi un halogène, un groupe alkyle en  $C_1-C_{60}$ , aryle en  $C_6-C_{60}$ , hétéroaryle en  $C_4-C_{60}$ , hétérocycloalkyle à 5 ou 6 éléments contenant un ou plusieurs hétéroatome(s) choisi(s) parmi N, O et S, cycloalkyle en  $C_3-C_{60}$ , tri(alkyle en  $C_1-C_{60}$ )Silyle, di(alkyle en  $C_1-C_{60}$ ) (aryle en  $C_6-C_{60}$ )silyle, tri(aryle en  $C_6-C_{60}$ ) silyle, adamantyle, bicycloalkyle en  $C_7-C_{60}$ , alcényle en  $C_2-C_{60}$ , alcynyle en  $C_2-C_{60}$ , cyano, (alkyle en  $C_1-C_{60}$ ) amino, (aryle en  $C_6-C_{60}$ )amino, (aryle en  $C_6-C_{60}$ ) (alkyle en  $C_1-C_{60}$ ), (alkyle en  $C_1-C_{60}$ )oxy, (alkyle en  $C_1-C_{60}$ )

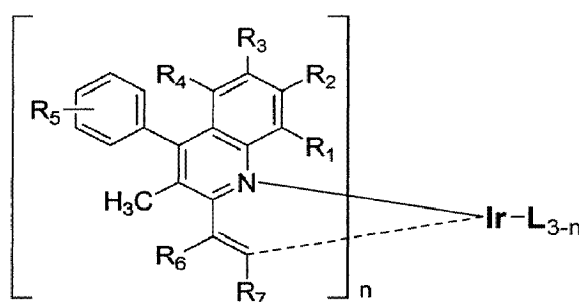
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thio, (aryle en C<sub>6</sub>-C<sub>60</sub>)oxy, (aryle en C<sub>6</sub>-C<sub>60</sub>)thio, (alcoxy en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, carboxyle, nitro et hydroxyle.

5. Dispositif électroluminescent organique selon la revendication 4, dans lequel la couche organique comprend un ou plusieurs composé(s) choisi(s) dans un groupe constitué de composés d'arylamine et de composés de styrylarylamine, ou d'un ou plusieurs métaux choisi(s) dans un groupe constitué de métaux organiques du groupe 1, du groupe 2, des métaux de transition de la quatrième période et de la cinquième période, des métaux de lanthanide et des éléments de transition d.

6. Dispositif électroluminescent organique selon la revendication 4, lequel est un affichage organique comprenant un composé électroluminescent organique représenté par un composé électroluminescent organique représenté par la formule chimique (1) :

formule chimique 1



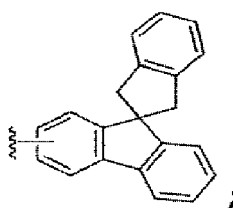
dans laquelle, L est un ligand organique ;

R<sub>1</sub> à R<sub>4</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, alcoxy en C<sub>1</sub>-C<sub>60</sub>, cycloalkyle en C<sub>3</sub>-C<sub>60</sub>, un halogène, un groupe tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle ou tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle ;

R<sub>5</sub> représente l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, un halogène ou un groupe aryle en C<sub>6</sub>-C<sub>60</sub> ;

R<sub>6</sub> et R<sub>7</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> avec ou sans substituant(s) d'halogène, aryle en C<sub>6</sub>-C<sub>60</sub>, un halogène, un groupe cyano, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, alcoxy en C<sub>1</sub>-C<sub>60</sub>, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino ou di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino, ou R<sub>6</sub> et R<sub>7</sub> peuvent être liés via un groupe alkylène en C<sub>3</sub>-C<sub>12</sub> ou alcénylène en C<sub>3</sub>-C<sub>12</sub> avec ou sans noyau condensé pour former un noyau alicyclique, ou un noyau aromatique monocyclique ou polycyclique ;

les groupes alkyle ou aryle de R<sub>6</sub> et R<sub>7</sub>, ou le noyau alicyclique, ou le noyau aromatique monocyclique ou polycyclique formé à partir de ceux-ci par liaison via un groupe alkylène en C<sub>3</sub>-C<sub>12</sub> ou alcénylène en C<sub>3</sub>-C<sub>12</sub> avec ou sans noyau condensé peuvent encore être substitués par un ou plusieurs substituant(s) choisi(s) parmi un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> avec ou sans substituant(s) d'halogène, un halogène, un groupe cyano, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, alcoxy en C<sub>1</sub>-C<sub>60</sub>, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino, di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino, phényle, naphtyle, anthryle, fluorényle, spiro-bifluorényle et

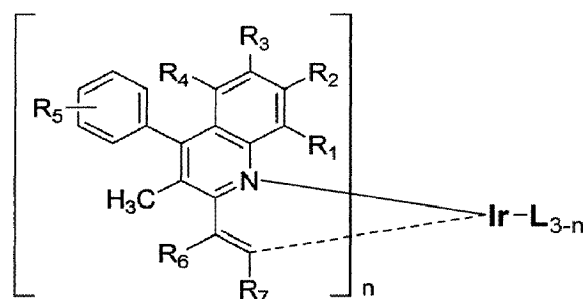


et

n est un nombre entier de 1 à 3 et des composés présentant le pic électroluminescent avec une longueur d'onde du bleu et du vert simultanément.

7. Dispositif électroluminescent organique selon la revendication 4, dans lequel la couche organique comprend une couche électroluminescente et une couche génératrice de charge.
8. Dispositif électroluminescent organique selon la revendication 4, dans lequel une région mixte d'un dopant réducteur et d'une substance organique, ou une région mixte d'un dopant oxydant et d'une substance organique est placée sur la surface interne d'une électrode ou des deux électrodes parmi la paire d'électrodes.
9. Pile solaire organique qui comprend un composé électroluminescent organique représenté par un composé électroluminescent organique représenté par la formule chimique (1) :

formule chimique 1



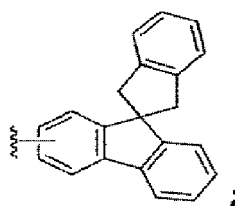
dans laquelle, L est un ligand organique ;

R<sub>1</sub> à R<sub>4</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, alcoxy en C<sub>1</sub>-C<sub>60</sub>, cycloalkyle en C<sub>3</sub>-C<sub>60</sub>, un halogène, un groupe tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle ou tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle ;

R<sub>5</sub> représente l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub>, un halogène ou un groupe aryle en C<sub>6</sub>-C<sub>60</sub> ;

R<sub>6</sub> et R<sub>7</sub> représentent indépendamment l'hydrogène, un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> avec ou sans substituant(s) d'halogène, un groupe aryle en C<sub>6</sub>-C<sub>60</sub>, un halogène, un groupe cyano, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, alcoxy en C<sub>1</sub>-C<sub>60</sub>, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino ou di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino, ou R<sub>6</sub> et R<sub>7</sub> peuvent être liés via un groupe alkylène en C<sub>3</sub>-C<sub>12</sub> ou alcénylène en C<sub>3</sub>-C<sub>12</sub> avec ou sans noyau condensé pour former un noyau alicyclique, ou un noyau aromatique monocyclique ou polycyclique ;

les groupes alkyle ou aryle de R<sub>6</sub> et R<sub>7</sub>, ou le noyau alicyclique, ou le noyau aromatique monocyclique ou polycyclique formé à partir de ceux-ci par liaison via un groupe alkylène en C<sub>3</sub>-C<sub>12</sub> ou alcénylène en C<sub>3</sub>-C<sub>12</sub> avec ou sans noyau condensé peuvent encore être substitués par un ou plusieurs substituant(s) choisi(s) parmi un groupe alkyle en C<sub>1</sub>-C<sub>60</sub> avec ou sans substituant(s) d'halogène, un halogène, un groupe cyano, tri(alkyle en C<sub>1</sub>-C<sub>60</sub>)silyle, tri(aryle en C<sub>6</sub>-C<sub>60</sub>)silyle, alcoxy en C<sub>1</sub>-C<sub>60</sub>, (alkyle en C<sub>1</sub>-C<sub>60</sub>)carbonyle, (aryle en C<sub>6</sub>-C<sub>60</sub>)carbonyle, di(alkyle en C<sub>1</sub>-C<sub>60</sub>)amino, di(aryle en C<sub>6</sub>-C<sub>60</sub>)amino, phényle, naphtyle, anthryle, fluorényle, spiro-bifluorényle et



et  
n est un nombre entier de 1 à 3.

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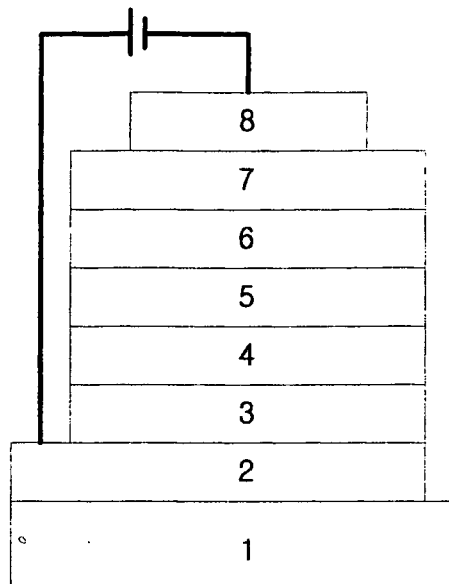
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Fig. 1



**REFERENCES CITED IN THE DESCRIPTION**

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**Non-patent literature cited in the description**

- **A. Tsuboyama et al.** *J. Am. Chem. Soc.*, 2003, vol. 125 (42), 12971-12979 [0004]

专利名称(译)	新型红色电致发光化合物和使用其的有机电致发光器件		
公开(公告)号	<a href="#">EP2066150B1</a>	公开(公告)日	2010-05-26
申请号	EP2008170051	申请日	2008-11-27
申请(专利权)人(译)	GRACEL显示增量.		
当前申请(专利权)人(译)	GRACEL显示增量.		
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摘要(译)

本发明涉及具有高发光效率的新型有机电致发光化合物，以及包含该化合物的有机电致发光器件。根据本发明的新型有机电致发光化合物由化学式(1)表示：

