



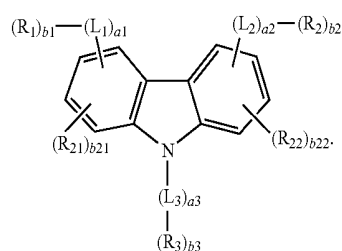
US 20160099424A1

(19) **United States**(12) **Patent Application Publication**
KIM et al.(10) **Pub. No.: US 2016/0099424 A1**(43) **Pub. Date: Apr. 7, 2016**(54) **CARBAZOLE-BASED COMPOUND AND
ORGANIC LIGHT-EMITTING DEVICE
INCLUDING THE SAME**(71) Applicant: **SAMSUNG DISPLAY CO., LTD.,**
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Seunggak YANG, Yongin-City (KR)(21) Appl. No.: **14/676,610**(22) Filed: **Apr. 1, 2015**(30) **Foreign Application Priority Data**

Oct. 1, 2014 (KR) 10-2014-0132491

Publication Classification(51) **Int. Cl.**
H01L 51/00 (2006.01)
C07D 487/04 (2006.01)
H01L 51/50 (2006.01)
C07D 471/04 (2006.01)(52) **U.S. CL.**
CPC **H01L 51/0072** (2013.01); **C07D 471/04**
(2013.01); **C07D 487/04** (2013.01); **H01L**
51/5016 (2013.01); **H01L 51/5012** (2013.01)(57) **ABSTRACT**

A carbazole-based compound and an organic light-emitting device, the carbazole-based compound being represented by Chemical Formula 1



<Formula 1>

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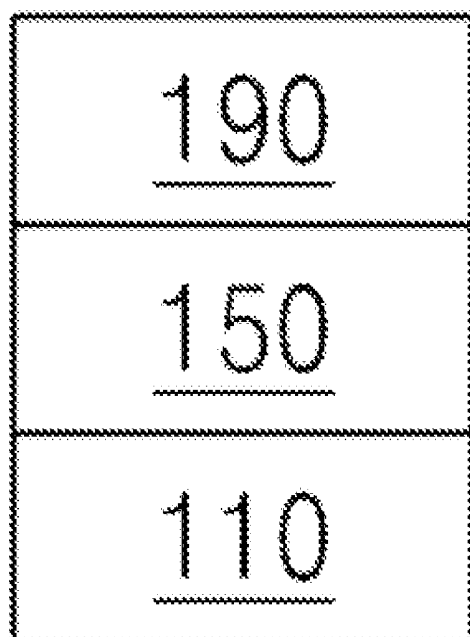
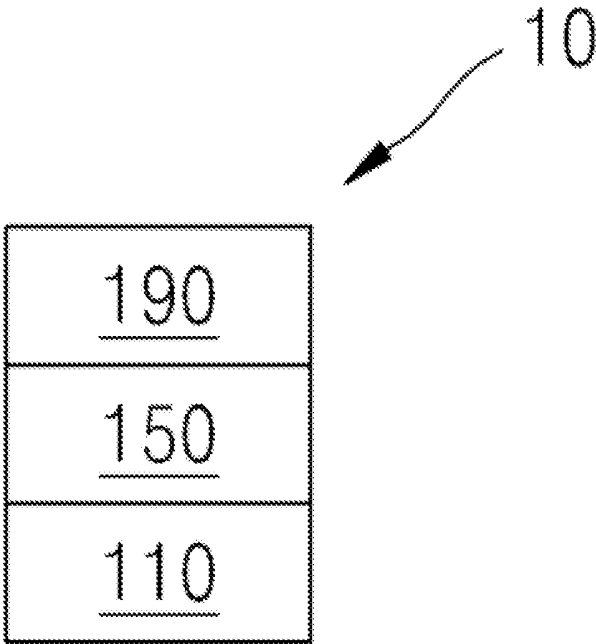


FIG. 1



CARBAZOLE-BASED COMPOUND AND ORGANIC LIGHT-EMITTING DEVICE INCLUDING THE SAME

CROSS-REFERENCE TO RELATED APPLICATION

[0001] Korean Patent Application No. 10-2014-0132491, filed on Oct. 1, 2014, in the Korean Intellectual Property Office, and entitled: "Carbazole-Based Compound and Organic Light-Emitting Device Including the Same," is incorporated by reference herein in its entirety.

BACKGROUND

[0002] 1. Field

[0003] Embodiments relate to a carbazole-based compound and an organic light-emitting device including the same.

[0004] 2. Description of the Related Art

[0005] Organic light emitting devices are self-emission devices that have wide viewing angles, high contrast ratios, short response times, and excellent brightness, driving voltage, and response speed characteristics, and produce full-color images.

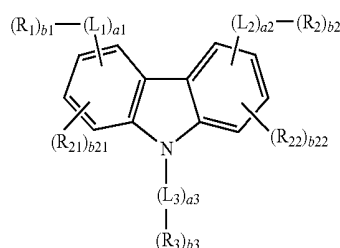
[0006] The organic light-emitting device may include a first electrode disposed on a substrate, and a hole transport region, an emission layer, an electron transport region, and a second electrode, which are sequentially disposed on the first electrode. Holes provided from the first electrode may move toward the emission layer through the hole transport region, and electrons provided from the second electrode may move toward the emission layer through the electron transport region. Carriers, such as holes and electrons, are recombined in the emission layer to produce excitons. These excitons change from an excited state to a ground state, thereby generating light.

SUMMARY

[0007] One or more exemplary embodiments include a novel carbazole-based compound and an organic light-emitting device including the same.

[0008] Additional aspects will be set forth in part in the description which follows and, in part, will be apparent from the description, or may be learned by practice of the presented embodiments.

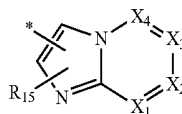
[0009] According to one or more exemplary embodiments, a carbazole-based compound may be represented by Formula 1 below:



<Formula 1>

-continued

<Formula 2>



[0010] wherein in the formulae above,

[0011] X_1 may be $C(R_{11})$ or N, X_2 may be $C(R_{12})$ or N, X_3 may be $C(R_{13})$ or N, and X_4 may be $C(R_{14})$ or N;

[0012] R_{11} to R_{15} , R_{21} and R_{22} may be each independently selected from a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a substituted or unsubstituted C_1 - C_{60} alkyl group, a substituted or unsubstituted C_2 - C_{60} alkenyl group, a substituted or unsubstituted C_2 - C_{60} alkynyl group, a substituted or unsubstituted C_1 - C_{60} alkoxy group, a substituted or unsubstituted C_3 - C_{10} cycloalkyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkyl group, a substituted or unsubstituted C_3 - C_{10} cycloalkenyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenyl group, a substituted or unsubstituted C_6 - C_{60} aryl group, a substituted or unsubstituted C_6 - C_{60} aryloxy group, a substituted or unsubstituted C_6 - C_{60} arylthio group, a substituted or unsubstituted C_1 - C_{60} heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, —Si(Q_1)(Q_2)(Q_3), —B(Q_4)(Q_5) and —N(Q_6)(Q_7);

[0013] b_{21} and b_{22} may be each independently selected from 0, 1, 2 and 3;

[0014] L_1 to L_3 may be each independently selected from a substituted or unsubstituted C_3 - C_{10} cycloalkylene group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkylene group, a substituted or unsubstituted C_3 - C_{10} cycloalkenylene group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenylene group, a substituted or unsubstituted C_6 - C_{60} arylene group, a substituted or unsubstituted C_1 - C_{60} heteroarylene group, a substituted or unsubstituted divalent non-aromatic condensed polycyclic group and a substituted or unsubstituted divalent non-aromatic condensed heteropolycyclic group;

[0015] a_1 to a_3 may be each independently selected from 0, 1, 2 and 3, and when a_1 is 2 or more, two or more of L_1 may be present and the two or more L_1 may be identical or different, and when a_2 is 2 or more, two or more L_2 may be present and the two or more L_2 may be identical or different, when a_3 is 2 or more, two or more L_3 may be present and two or more L_3 may be identical or different;

[0016] R_1 may be a substituted or unsubstituted carbazolyl group;

[0017] R_2 and R_3 may be each independently selected from a group represented by Formula 2, a hydrogen, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a substituted or unsubstituted C_1 - C_{60} alkyl group, a substituted or unsubstituted C_2 - C_{60} alkenyl group, a substituted or unsubstituted C_2 - C_{60} alkynyl group, a substituted or unsubstituted C_1 - C_{60} alkoxy group, a substituted or unsubstituted C_3 - C_{10} cycloalkyl group, a substituted or unsubstituted C_1 - C_{10} het-

erocycloalkyl group, a substituted or unsubstituted C₃-C₁₀ cycloalkenyl group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkenyl group, a substituted or unsubstituted C₆-C₆₀ aryl group, a substituted or unsubstituted C₁-C₆₀ heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group (wherein the “substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group” does not include the group represented by Formula 2 above), —Si(Q₁)(Q₂)(Q₃), —B(Q₄)(Q₅) and —N(Q₆)(Q₇);

[0018] at least one of R₂ and R₃ is a group represented by Formula 2;

[0019] b1 to b3 may be each independently an integer of 1 to 3;

[0020] at least one of substituents of the substituted carbazolyl group, substituted C₃-C₁₀ cycloalkylene group, substituted C₁-C₁₀ heterocycloalkylene group, substituted C₃-C₁₀ cycloalkenylene group, substituted C₁-C₁₀ heterocycloalkenylene group, substituted C₆-C₆₀ arylene group, substituted C₁-C₆₀ heteroarylene group, substituted divalent non-aromatic condensed polycyclic group, substituted divalent non-aromatic condensed heteropolycyclic group, substituted C₁-C₆₀ alkyl group, substituted C₂-C₆₀ alkenyl group, substituted C₂-C₆₀ alkynyl group, substituted C₁-C₆₀ alkoxy group, substituted C₃-C₁₀ cycloalkyl group, substituted C₁-C₁₀ heterocycloalkyl group, substituted C₃-C₁₀ cycloalkenyl group, substituted C₁-C₁₀ heterocycloalkenyl group, substituted C₆-C₆₀ aryl group, substituted C₆-C₆₀ aryloxy group, substituted C₆-C₆₀ arylthio group, substituted C₁-C₆₀ heteroaryl group, substituted monovalent non-aromatic condensed polycyclic group, and substituted monovalent non-aromatic condensed heteropolycyclic group may be selected from

[0021] a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group;

[0022] a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group and a C₁-C₆₀ alkoxy group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, —Si(Q₁₁)(Q₁₂)(Q₁₃), —B(Q₁₄)(Q₁₅) and —N(Q₁₆)(Q₁₇);

[0023] a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic heterocondensed polycyclic group;

[0024] a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ het-

erocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group and a monovalent non-aromatic condensed heteropolycyclic group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, —Si(Q₂₁)(Q₂₂)(Q₂₃), —B(Q₂₄)(Q₂₅) and —N(Q₂₆)(Q₂₇); and

[0025] —Si(Q₃₁)(Q₃₂)(Q₃₃), —B(Q₃₄)(Q₃₅) and —N(Q₃₆)(Q₃₇);

[0026] wherein Q₁ to Q₇, Q₁₁ to Q₁₇, Q₂₁ to Q₂₇, and Q₃₁ to Q₃₇ may be each independently selected from a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group.

[0027] Another aspect provides an organic light-emitting device including: a first electrode; a second electrode facing the first electrode; and an organic layer that is disposed between the first electrode and the second electrode and includes an emission layer, wherein the organic layer includes at least one of the carbazole-based compounds described above.

BRIEF DESCRIPTION OF THE DRAWING

[0028] Features will be apparent to those of skill in the art by describing in detail exemplary embodiments with reference to the attached drawings in which:

[0029] FIG. 1 illustrates a schematic view of an organic light-emitting device according to an embodiment.

DETAILED DESCRIPTION

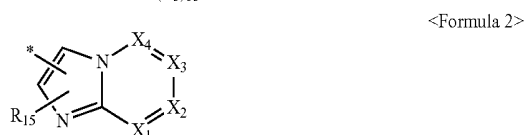
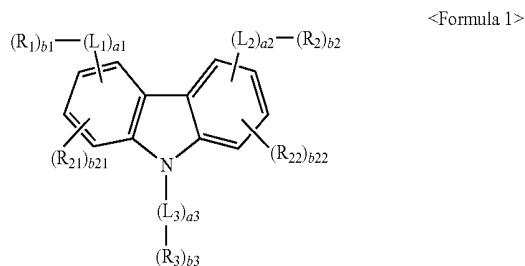
[0030] Example embodiments will now be described more fully hereinafter with reference to the accompanying drawing; however, they may be embodied in different forms and should not be construed as limited to the embodiments set forth herein. Rather, these embodiments are provided so that this disclosure will be thorough and complete, and will fully convey exemplary implementations to those skilled in the art.

[0031] In the drawing FIGURE, the dimensions of layers and regions may be exaggerated for clarity of illustration. Like reference numerals refer to like elements throughout.

[0032] As used herein, the term “and/or” includes any and all combinations of one or more of the associated listed items. Expressions such as “at least one of” when preceding a list of

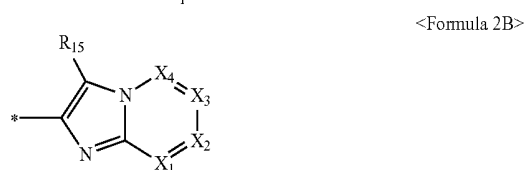
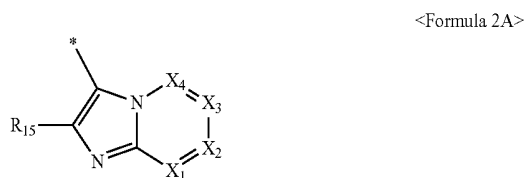
elements, modify the entire list of elements and do not modify the individual elements of the list.

[0033] A carbazole-based compound may be represented by Formula 1 below, and at least one of R_2 and R_3 in Formula 1 is represented by Formula 2 below.



[0034] b_2 and b_3 in Formula 1 are not 0, and thus the carbazole-based compound necessarily includes a group represented by Formula 2 as a substituent.

[0035] Formula 2 may be represented by Formula 2A or 2B below.



[0036] In Formula 1, X_1 may be $C(R_{11})$ or N, X_2 may be $C(R_{12})$ or N, X_3 may be $C(R_{13})$ or N, and X_4 may be $C(R_{14})$ or N.

[0037] In some embodiments, in the formulae, X_1 may be $C(R_{11})$, X_2 may be $C(R_{12})$, X_3 may be $C(R_{13})$, and X_4 may be $C(R_{14})$.

[0038] In some embodiments, X_1 may be N, X_2 may be $C(R_{12})$, X_3 may be $C(R_{13})$, and X_4 may be $C(R_{14})$, but are not limited thereto.

[0039] In the formulae above, R_{11} to R_{15} , R_{21} and R_{22} may be each independently selected from or include, e.g., a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a substituted or unsubstituted C_1 - C_{60} alkyl group, a substituted or unsubstituted C_2 - C_{60} alkenyl group, a substituted or unsubstituted C_2 - C_{60} alkynyl group, a substituted or unsubstituted C_1 - C_{60} alkoxy group, a substituted or unsubstituted C_3 - C_{10} cycloalkyl

group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkyl group, a substituted or unsubstituted C_3 - C_{10} cycloalkenyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenyl group, a substituted or unsubstituted C_6 - C_{60} aryl group, a substituted or unsubstituted C_6 - C_{60} aryloxy group, a substituted or unsubstituted C_6 - C_{60} arylthio group, a substituted or unsubstituted C_1 - C_{60} heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, —Si(Q_1)(Q_2)(Q_3), —B(Q_4)(Q_5) and —N(Q_6)(Q_7).

[0040] According to an embodiment, in the formulae above, R_{11} to R_{15} , R_{21} and R_{22} may be each independently selected from:

[0041] a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, and a C_1 - C_{20} alkoxy group;

[0042] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group and a thiadiazolyl group; and

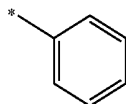
[0043] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group and a thiadiazolyl group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, C_1 - C_{20} alkyl group, C_1 - C_{20} alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl

group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazoyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group and a thiadiazolyl group.

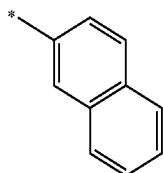
[0044] According to another embodiment, in the formulae above R_{11} to R_{15} , R_{21} and R_{22} may be each independently selected from:

[0045] a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, C_1 - C_{20} alkyl group and C_1 - C_{20} alkoxy group; and

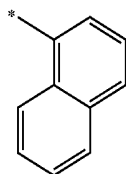
[0046] a group represented by one of Formulae 7-1 to 7-45 below, but are not limited thereto:



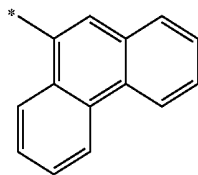
Formula 7-1



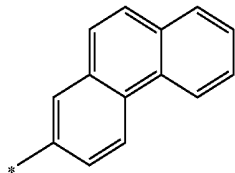
Formula 7-2



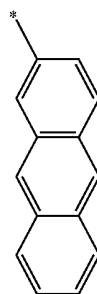
Formula 7-3



Formula 7-4

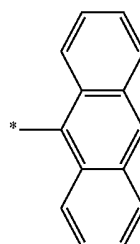


Formula 7-5

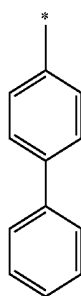


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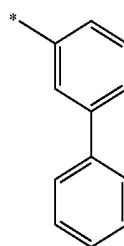
Formula 7-6



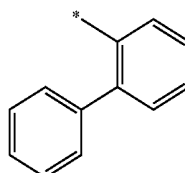
Formula 7-7



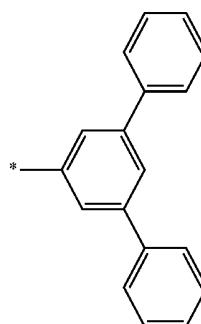
Formula 7-8



Formula 7-9

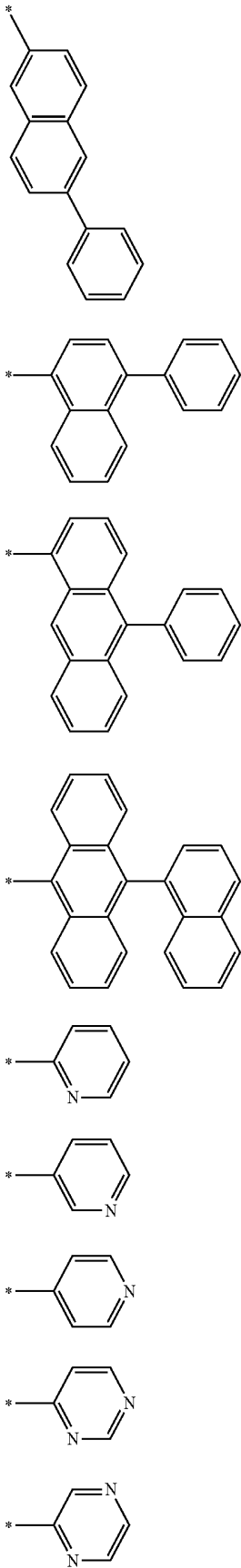


Formula 7-10



Formula 7-11

-continued



Formula 7-12

Formula 7-13

Formula 7-14

Formula 7-15

Formula 7-16

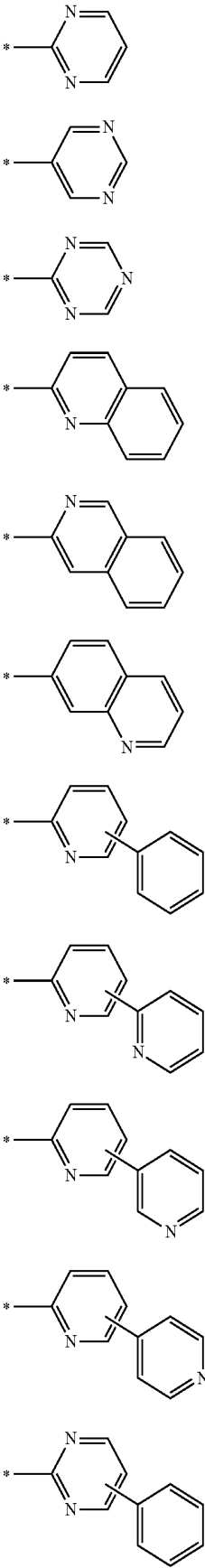
Formula 7-17

Formula 7-18

Formula 7-19

Formula 7-20

-continued



Formula 7-21

Formula 7-22

Formula 7-23

Formula 7-24

Formula 7-25

Formula 7-26

Formula 7-27

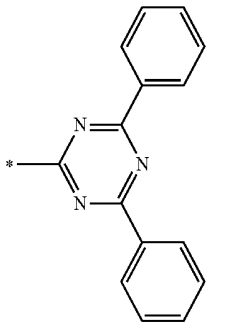
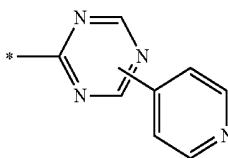
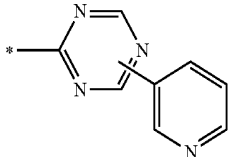
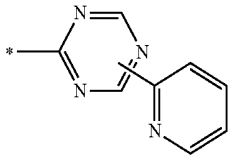
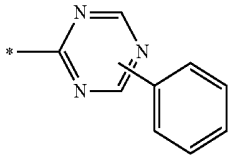
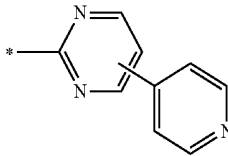
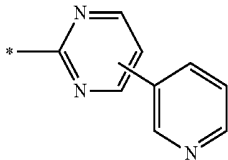
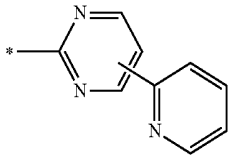
Formula 7-28

Formula 7-29

Formula 7-30

Formula 7-31

-continued



Formula 7-32

Formula 7-33

Formula 7-34

Formula 7-35

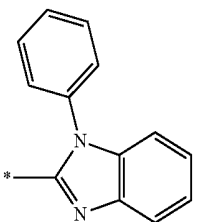
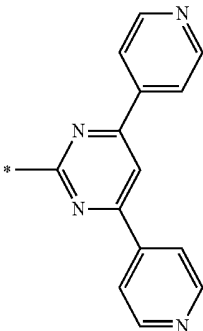
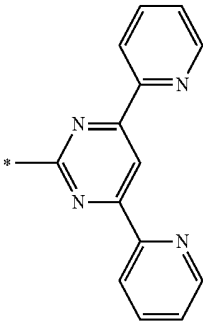
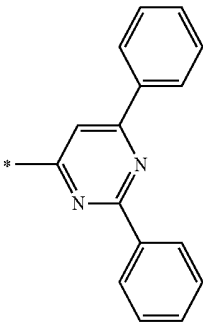
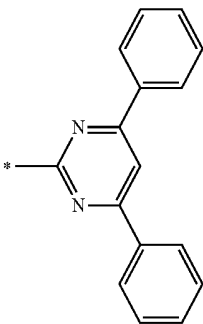
Formula 7-36

Formula 7-37

Formula 7-38

Formula 7-39

-continued



Formula 7-40

Formula 7-41

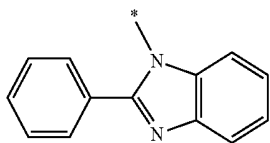
Formula 7-42

Formula 7-43

Formula 7-44

-continued

Formula 7-45



[0047] * in Formulae 7-1 to 7-45 indicates a binding site to a neighboring atom.

[0048] In some embodiments, in the formulae above,

[0049] R_{11} to R_{15} may be each independently selected from:

[0050] a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group and a C_1 - C_{20} alkoxy group; and

[0051] a group represented by one of Formulae 7-1 to 7-45 above; and

[0052] R_{21} and R_{22} may be each independently selected from:

[0053] a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, and a C_1 - C_{20} alkoxy group.

[0054] b21 in Formula 1 indicates the number of R_{21} , and b21 may be selected from 0, 1, 2 and 3. When b21 is 2 or more, two or more R_{21} may be present, and the two or more R_{21} may be identical or different. b22 in Formula 1 may be understood based on the description of b21 defined herein.

[0055] L_1 to L_3 in Formula 1 may be each independently selected from or include, e.g., a substituted or unsubstituted C_3 - C_{10} cycloalkylene group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkylene group, a substituted or unsubstituted C_3 - C_{10} cycloalkenylene group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenylene group, a substituted or unsubstituted C_6 - C_{60} arylene group, a substituted or unsubstituted C_1 - C_{60} heteroarylene group, a substituted or unsubstituted divalent non-aromatic condensed polycyclic group and a substituted or unsubstituted divalent non-aromatic condensed heteropolycyclic group.

[0056] According to an embodiment, L_1 to L_3 in Formula 1 may be each independently selected from:

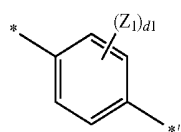
[0057] a phenylene group, a pentalenylene group, an indenylene group, a naphthylene group, an azulenylene group, a heptalenylene group, an indacenylene group, an acenaphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzofluorenylene group, a dibenzofluorenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylene group, a fluoranthenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a naphthacenylene group, a picenylene group, a perylenylene group, a pentaphenylene group, a hexacenylene group, a pentacenylene group, a rubicenylene group, a coronenylene group, an ovalenylene group, a pyrrolylene group, a thiophenylene group, a furanylene group, an imidazolylene group, a pyrazolylene group, a thiazolylene group, an isothiazolylene group, an oxazolylene group, a isooxazolylene group, a pyridinylene group, a pyrazinylene group, a pyrimidinylene group, a pyridazinylene group, an isoidolylene

group, an indolylene group, an indazolylene group, a purinylene group, a quinolinylenylene group, an isoquinolinylenylene group, a benzoquinolinylenylene group, a phthalazinylene group, a naphthyridinylene group, a quinoxalinylene group, a quinazolinylene group, a cinnolinylenylene group, a carbazolylene group, a phenanthridinylene group, an acridinylene group, a phenanthrolinylene group, a phenazinylene group, a benzoimidazolylene group, a benzofuranylenylene group, a benzothienophenylene group, an isobenzothiazolylene group, a benzooxazolylene group, an isobenzooxazolylene group, a triazolylene group, a tetrazolylene group, an oxadiazolylene group, a triazinylene group, a dibenzofuranylenylene group, a dibenzothiophenylene group, a benzocarbazolylene group, a dibenzocarbazolylene group, a thiadiazolylene group, an imidazopyridinylene group and an imidazopyrimidinylene group; and

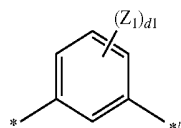
[0058] a phenylene group, a pentalenylene group, an indenylene group, a naphthylene group, an azulenylene group, a heptalenylene group, an indacenylene group, an acenaphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzofluorenylene group, a dibenzofluorenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylene group, a fluoranthenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a naphthacenylene group, a picenylene group, a perylenylene group, a pentaphenylene group, a hexacenylene group, a pentacenylene group, a rubicenylene group, a coronenylene group, ovalenylene group, a pyrrolylene group, a thiophenylene group, a furanylene group, an imidazolylene group, a pyrazolylene group, a thiazolylene group, an isothiazolylene group, an oxazolylene group, a isooxazolylene group, a pyridinylene group, a pyrazinylene group, a pyrimidinylene group, a pyridazinylene group, an isoidolylene group, an indolylene group, an indazolylene group, a purinylene group, a quinolinylenylene group, an isoquinolinylenylene group, a benzoquinolinylenylene group, a phthalazinylene group, a naphthyridinylene group, a quinoxalinylene group, a quinazolinylene group, a cinnolinylenylene group, a carbazolylene group, a phenanthridinylene group, an acridinylene group, a phenanthrolinylene group, a phenazinylene group, a benzoimidazolylene group, a benzofuranylenylene group, a benzothiophenylene group, an isobenzothiazolylene group, a benzooxazolylene group, an isobenzooxazolylene group, a triazolylene group, a tetrazolylene group, an oxadiazolylene group, a triazinylene group, a dibenzofuranylenylene group, a dibenzothiophenylene group, a benzocarbazolylene group, a dibenzocarbazolylene group, a thiadiazolylene group, an imidazopyridinylene group and an imidazopyrimidinylene group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, C_1 - C_{20} alkyl group, C_1 - C_{20} alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenyl group, a heptalenyl group, an indacenyl group, an acenaphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl

group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, a tetrazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzo-carbazolyl group, a dibenzocarbazolyl group, a thiadiazolyl group, an imidazopyridinylene group and an imidazopyrimidinylene group.

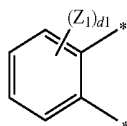
[0059] According to another embodiment, L_1 to L_3 in Formula 1 may be each independently selected from a group represented by one of Formulae 3-1 to 3-33 below, but are not limited thereto:



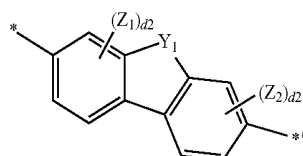
Formula 3-1



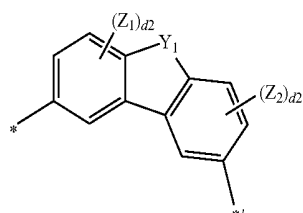
Formula 3-2



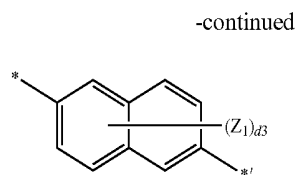
Formula 3-3



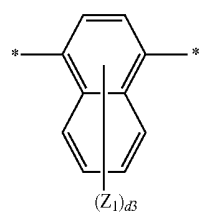
Formula 3-4



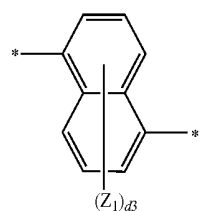
Formula 3-5



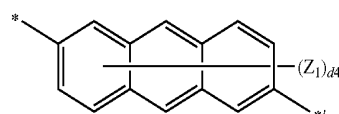
Formula 3-6



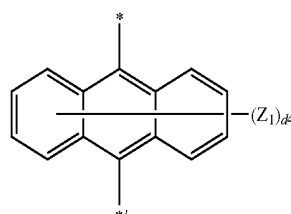
Formula 3-7



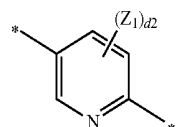
Formula 3-8



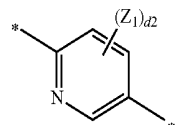
Formula 3-9



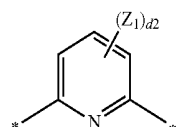
Formula 3-10



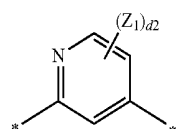
Formula 3-11



Formula 3-12

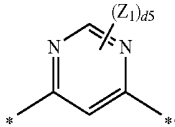
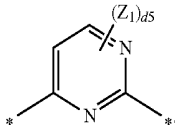
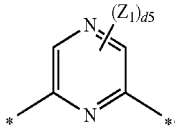
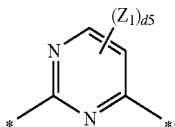
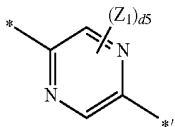
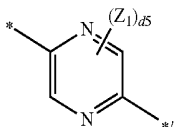
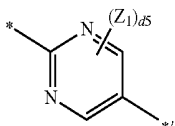
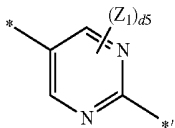
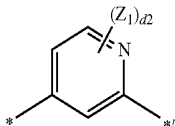
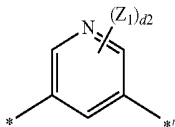


Formula 3-13



Formula 3-14

-continued



Formula 3-15

Formula 3-16

Formula 3-17

Formula 3-18

Formula 3-19

Formula 3-20

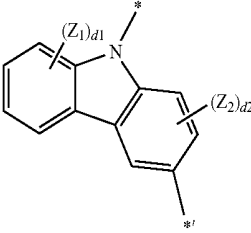
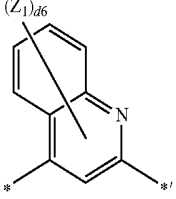
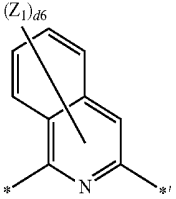
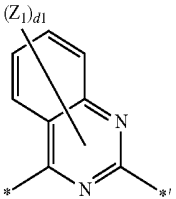
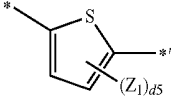
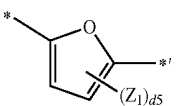
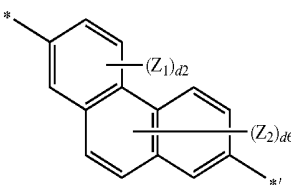
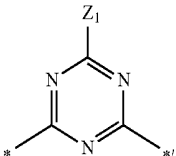
Formula 3-21

Formula 3-22

Formula 3-23

Formula 3-24

-continued



Formula 3-25

Formula 3-26

Formula 3-27

Formula 3-28

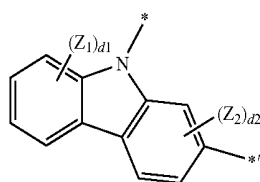
Formula 3-29

Formula 3-30

Formula 3-31

Formula 3-32

-continued



Formula 3-33

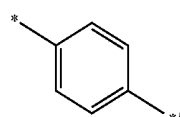
[0060] Wherein, in Formulae 3-1 to 3-33,

[0061] Y_1 may be O, S, C(Z_3)(Z_4), N(Z_5), or Si(Z_6)(Z_7);

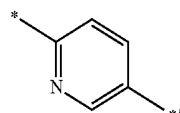
[0062] Z_1 to Z_7 may be each independently selected from a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 to C_{20} alkyl group, a C_1 to C_{20} alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, and a triazinyl group;

[0063] d_1 may be an integer of 1 to 4, d_2 may be an integer of 1 to 3, d_3 may be an integer of 1 to 6, d_4 may be an integer of 1 to 8, d_5 may be 1 or 2, and d_6 may be an integer of 1 to 5, and * and *' indicate binding sites to a neighboring atom.

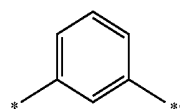
[0064] According to another embodiment, L_1 to L_3 in Formula 1 may be each independently selected from a group represented by one of Formulae 4-1 to 4-24 below, but are not limited thereto.



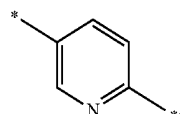
Formula 4-1



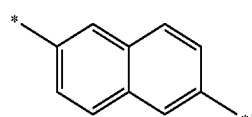
Formula 4-2



Formula 4-3

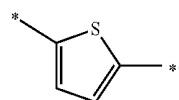


Formula 4-4

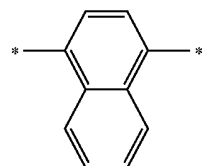


Formula 4-5

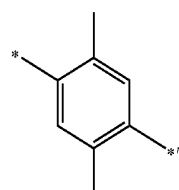
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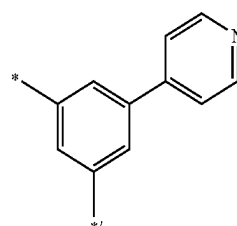
Formula 4-6



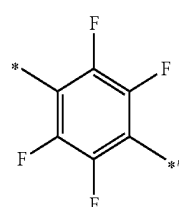
Formula 4-7



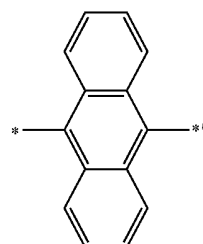
Formula 4-8



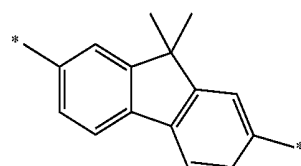
Formula 4-9



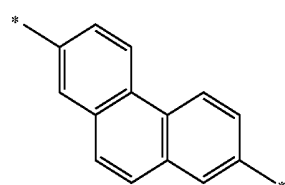
Formula 4-10



Formula 4-11

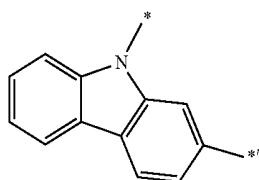
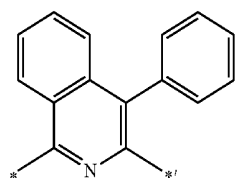
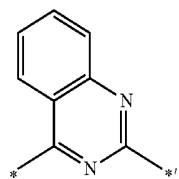
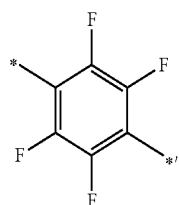
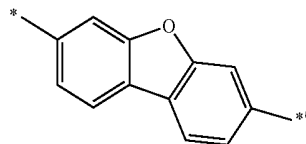
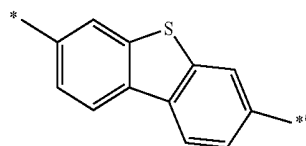
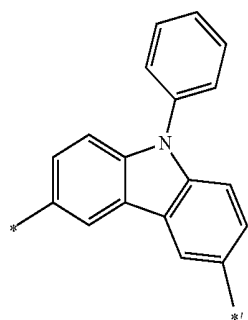


Formula 4-12



Formula 4-13

-continued



Formula 4-14

Formula 4-15

Formula 4-16

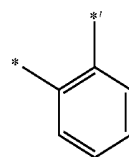
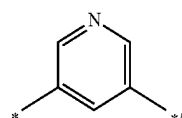
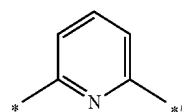
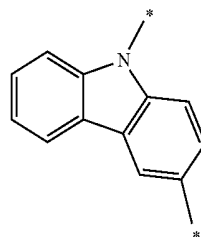
Formula 4-17

Formula 4-18

Formula 4-19

Formula 4-20

-continued



Formula 4-21

Formula 4-22

Formula 4-23

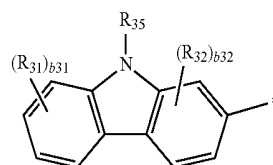
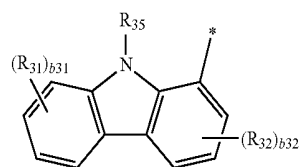
Formula 4-24

[0065] * and *' in Formulae 4-1 to 4-24 indicate binding sites to a neighboring atom.

[0066] a₁ in Formula 1 indicates the number of L₁, and a₁ may be selected from 0, 1, 2 and 3. When a₁ is 0, *-(L₁)_{a1}-* is a single bond. When a₁ is 2 or more, two or more L₁ may be present and the two or more L₁ may be identical or different. a₁ in Formula 1 may be 0 or 1. a₂ and a₃ in Formula 1 may be understood based on the description of a₁ defined herein.

[0067] In an implementation, R₁ in Formula 1 may be a substituted or unsubstituted carbazolyl group.

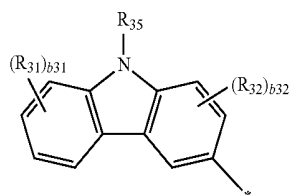
[0068] According to an embodiment, R₁ in Formula 1 may be selected from a group represented by one of Formulae 8-1 to 8-5 below.



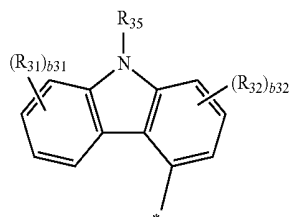
<Formula 8-1>

<Formula 8-2>

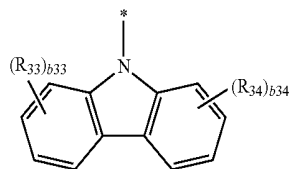
-continued



<Formula 8-3>



<Formula 8-4>



<Formula 8-5>

[0069] In Formulae 8-1 to 8-5,

[0070] R_{31} to R_{34} may be the same as R_{11} defined herein;

[0071] R_{35} may be selected from or include, e.g., a substituted or unsubstituted C_1 - C_{60} alkyl group, a substituted or unsubstituted C_2 - C_{60} alkenyl group, a substituted or unsubstituted C_2 - C_{60} alkynyl group, a substituted or unsubstituted C_1 - C_{60} alkoxy group, a substituted or unsubstituted C_3 - C_{10} cycloalkyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkyl group, a substituted or unsubstituted C_3 - C_{10} cycloalkenyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenyl group, a substituted or unsubstituted C_6 - C_{60} aryl group, a substituted or unsubstituted C_1 - C_{60} heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, $-\text{Si}(Q_1)(Q_2)(Q_3)$, $-\text{B}(Q_4)(Q_5)$ and $-\text{N}(Q_6)(Q_7)$ (Q_1 to Q_7 are already described above);

[0072] b_{31} , b_{33} , and b_{34} may be each independently selected from 0, 1, 2, 3, and 4, and b_{32} may be selected from 0, 1, 2, and 3.

[0073] For example, R_1 in Formula 1 may be represented by Formula 8-3 or 8-5.

[0074] R_2 and R_3 in Formula 1 may be each independently selected from or include, e.g., a group represented by Formula 2, a hydrogen, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a substituted or unsubstituted C_1 - C_{60} alkyl group, a substituted or unsubstituted C_2 - C_{60} alkenyl group, a substituted or unsubstituted C_2 - C_{60} alkynyl group, a substituted or unsubstituted C_1 - C_{60} alkoxy group, a substituted or unsubstituted C_3 - C_{10} cycloalkyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkyl group, a substituted or unsubstituted C_3 - C_{10} cycloalkenyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenyl group, a substituted or unsubstituted C_6 - C_{60} aryl group, a substituted or unsubstituted C_1 - C_{60} heteroaryl group, a substituted or unsubstituted monovalent non-

aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group (wherein the “substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group” does not include the group represented by Formula 2 above), $-\text{Si}(Q_1)(Q_2)(Q_3)$, $-\text{B}(Q_4)(Q_5)$ and $-\text{N}(Q_6)(Q_7)$, and at least one of R_2 and R_3 may be a group represented by Formula 2.

[0075] According to an embodiment, R_2 in Formula 1 may be selected from:

[0076] a group represented by Formula 2;

[0077] a hydrogen, a deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, and a C_1 - C_{20} alkoxy group;

[0078] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranlyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group; and

[0079] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranlyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group, each substituted with at least one selected from a deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, C_1 - C_{20} alkyl group, C_1 - C_{20} alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an

oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazoyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group.

[0080] R_3 may be selected from:

[0081] a group represented by Formula 2;

[0082] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazoyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group; and

[0083] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazoyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group,

a carbazoyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group.

[0084] At least one of R_2 and R_3 may be a group represented by Formula 2.

[0085] According to another embodiment, in Formula 1,

[0086] R_2 may be selected from:

[0087] a group represented by Formula 2;

[0088] a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group and a C_1 - C_{20} alkoxy group; and

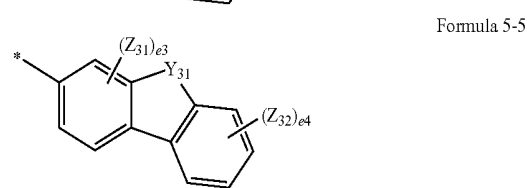
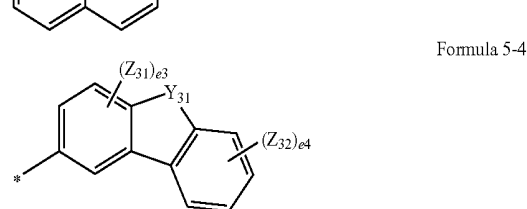
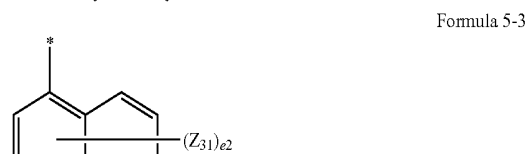
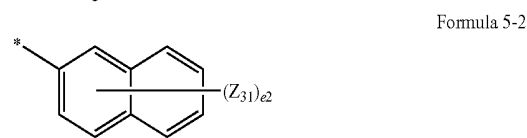
[0089] a group represented by one of Formulae 5-1 to 5-26 below.

[0090] R_3 may be selected from:

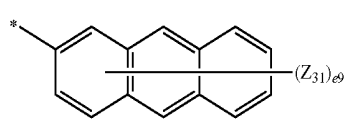
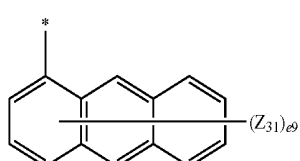
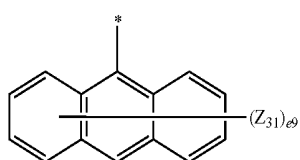
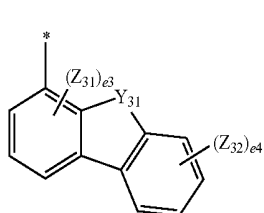
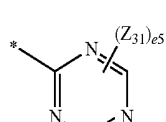
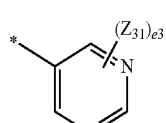
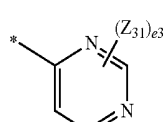
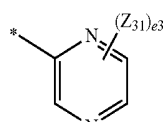
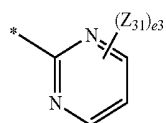
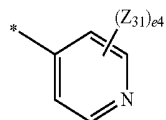
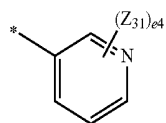
[0091] a group represented by Formula 2; and

[0092] a group represented by one of Formulae 5-1 to 5-26 below.

[0093] At least one of R_2 and R_3 may be a group represented by Formula 2, but is not limited thereto.



-continued



Formula 5-7

Formula 5-8

Formula 5-9

Formula 5-10

Formula 5-11

Formula 5-12

Formula 5-13

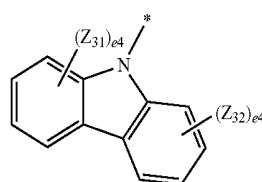
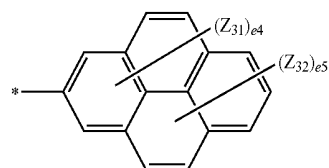
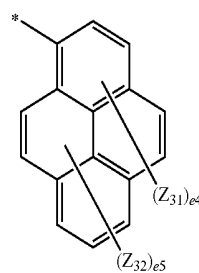
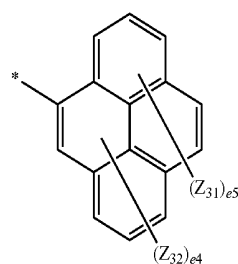
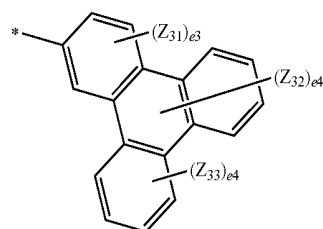
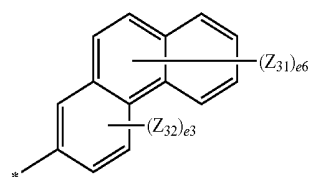
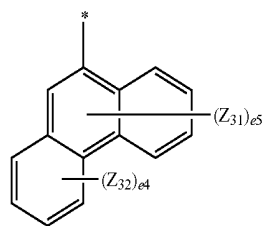
Formula 5-14

Formula 5-15

Formula 5-16

Formula 5-17

-continued



Formula 5-18

Formula 5-19

Formula 5-20

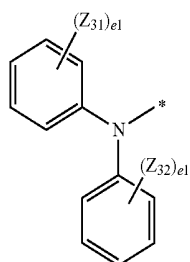
Formula 5-21

Formula 5-22

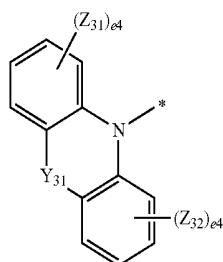
Formula 5-23

Formula 5-24

-continued



Formula 5-25



Formula 5-26

[0094] In Formulae 5-1 to 5-26,

[0095] Y_{31} may be O, S, $C(Z_{34})(Z_{35})$, $N(Z_{36})$, or $Si(Z_{37})(Z_{38})$;

[0096] Z_{31} to Z_{38} may be each independently selected from:

[0097] a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, and a C_1 - C_{20} alkoxy group;

[0098] a C_1 - C_{20} alkyl group and a C_1 - C_{20} alkoxy group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof;

[0099] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, and a dibenzocarbazolyl group;

[0100] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, and a dibenzocarbazolyl group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof,

thereof, a C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a phenyl group, and a naphthyl group; and

[0101] — $Si(Q_{31})(Q_{32})(Q_{33})$,

[0102] wherein Q_{31} to Q_{33} may be each independently selected from a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzo-fluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, and a triazinyl group.

[0103] e_1 may be an integer of 1 to 5, e_2 may be an integer of 1 to 7, e_3 may be an integer of 1 to 3, e_4 may be an integer of 1 to 4, e_5 may be 1 or 2, e_6 may be an integer of 1 to 6, e_9 may be an integer of 1 to 9, and * indicates a binding site to a neighboring atom.

[0104] According to another embodiment, in Formula 1,

[0105] R_2 may be selected from:

[0106] a group represented by Formula 2;

[0107] a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, and a C_1 - C_{20} alkoxy group; and

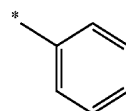
[0108] a group represented by one of Formulae 6-1 to 6-35 below.

[0109] R_3 may be selected from:

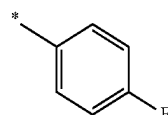
[0110] a group represented by Formula 2; and

[0111] a group represented by one of Formulae 6-1 to 6-35 below.

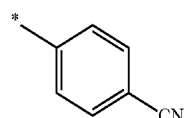
[0112] At least one of R_2 and R_3 may be a group represented by Formula 2, but is not limited thereto.



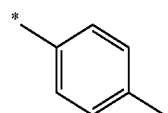
Formula 6-1



Formula 6-2

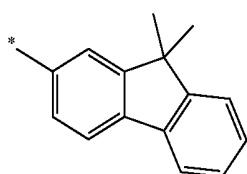
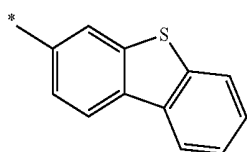
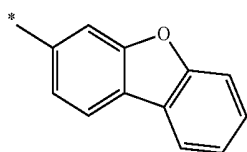
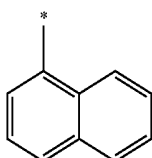
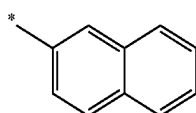
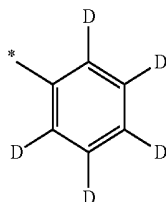
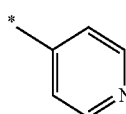
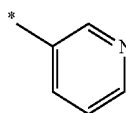
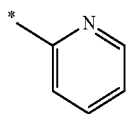


Formula 6-3



Formula 6-4

-continued



Formula 6-5

Formula 6-6

Formula 6-7

Formula 6-8

Formula 6-9

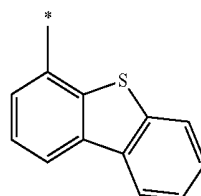
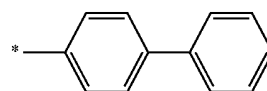
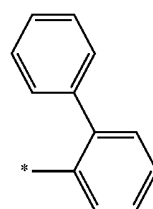
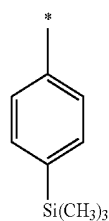
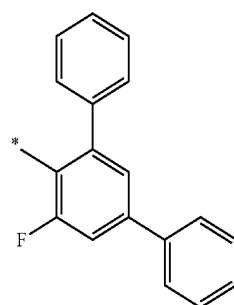
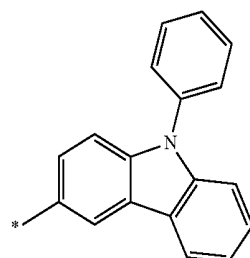
Formula 6-10

Formula 6-11

Formula 6-12

Formula 6-13

-continued



Formula 6-14

Formula 6-15

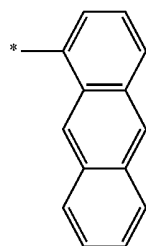
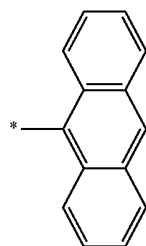
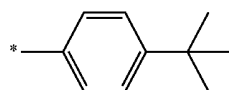
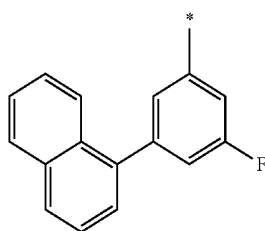
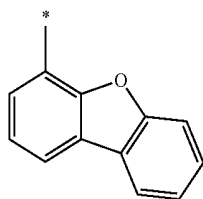
Formula 6-16

Formula 6-17

Formula 6-18

Formula 6-19

-continued



Formula 6-20

Formula 6-21

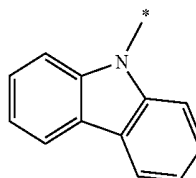
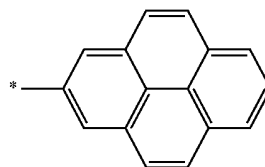
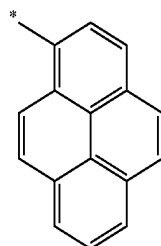
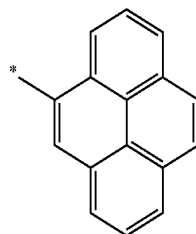
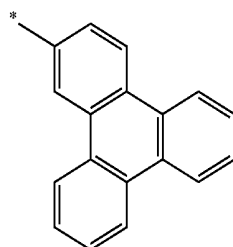
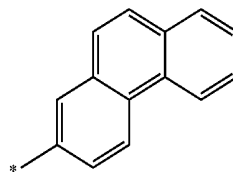
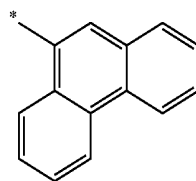
Formula 6-22

Formula 6-23

Formula 6-24

Formula 6-25

-continued



Formula 6-26

Formula 6-27

Formula 6-28

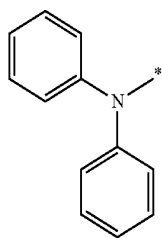
Formula 6-29

Formula 6-30

Formula 6-31

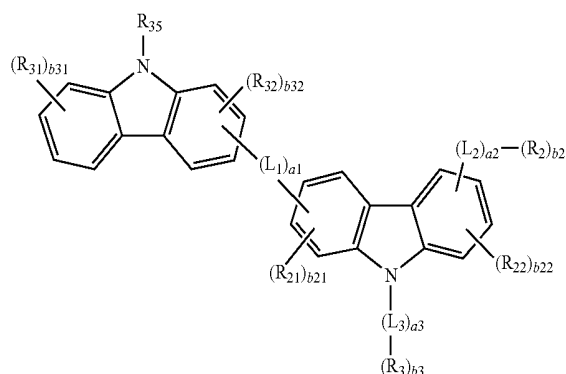
Formula 6-32

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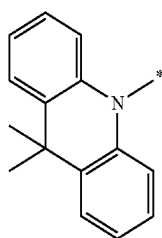
Formula 6-33

<Formula 1A>

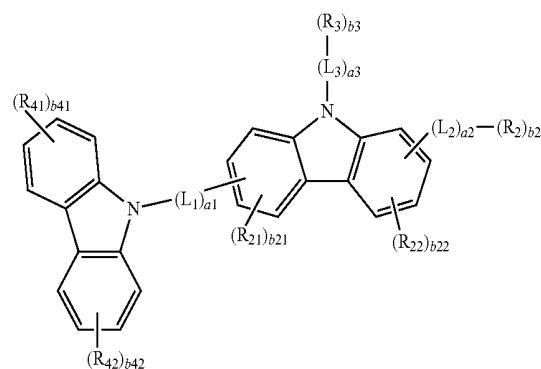
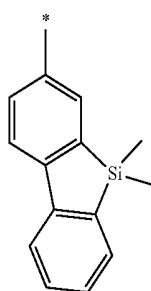


Formula 6-34

<Formula 1B>



Formula 6-35



[0113] * in Formulae 6-1 to 6-35 indicates a binding site to a neighboring atom.

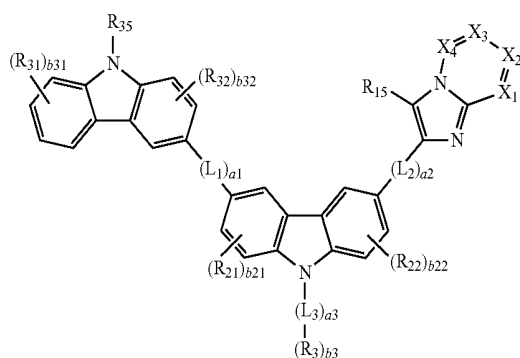
[0114] b1 in Formula 1 indicates the number of R_1 , and b1 may be integer of 1 to 3. When b1 is 2 or more, two or more R_1 may be present and the two or more R_1 may be identical or different. b2 and b3 may be understood based on the description of b1 defined herein.

[0115] According to an embodiment, a carbazole-based compound represented by Formula 1 may be represented by Formula 1A or Formula 1B below.

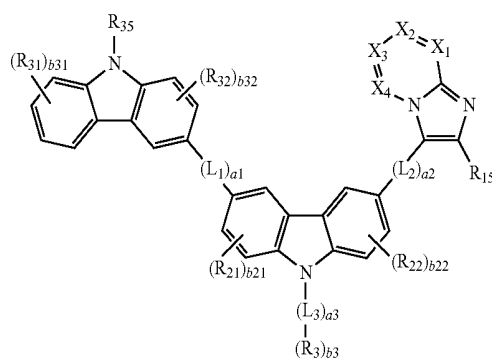
[0116] In Formulae 1A and 1B, L_1 to L_3 , a1 to a3, R_2 , R_3 , b2, b3, R_{21} , R_{22} , R_{35} , b21 and b22 may be the same as defined herein; R_{31} , R_{32} , R_{41} and R_{42} may be the same as R_{11} defined herein; b31, b41 and b42 may be each independently selected from 0, 1, 2, 3 and 4; and b32 may be selected from 0, 1, 2, and 3.

[0117] According to another embodiment, a carbazole-based compound represented by Formula 1 may be represented by one selected from Formulae 1A(1) to 1A(6) and Formulae 1B(1) to 1B(6):

<Formula 1A(1)>

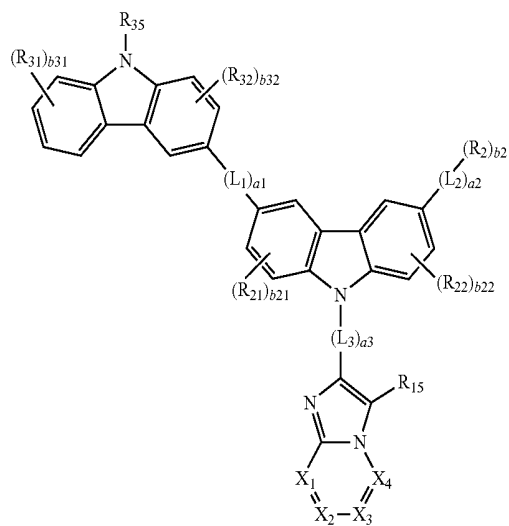


<Formula 1A(2)>

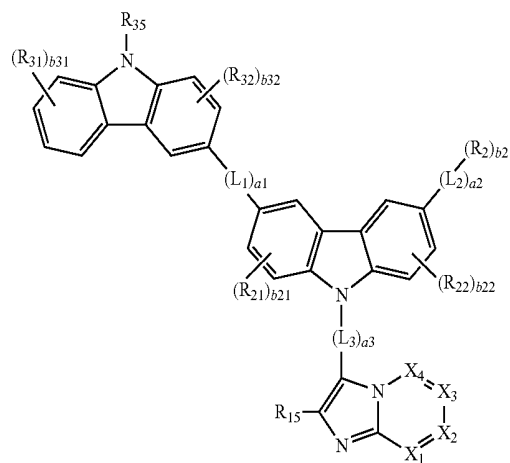


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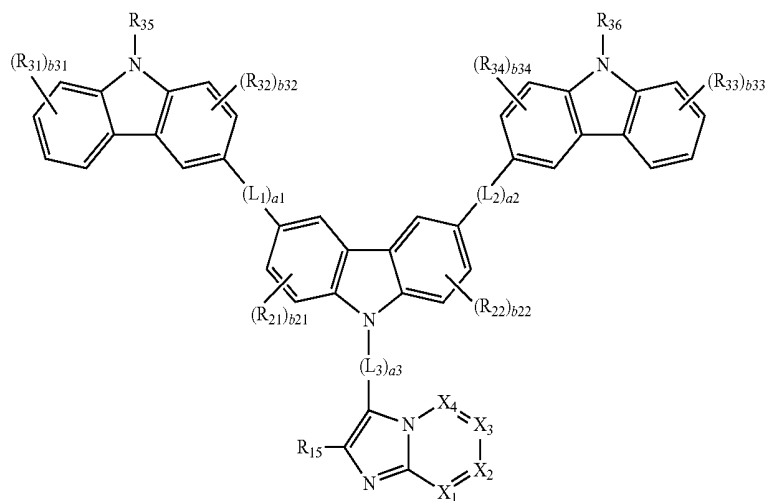
<Formula 1A(3)>



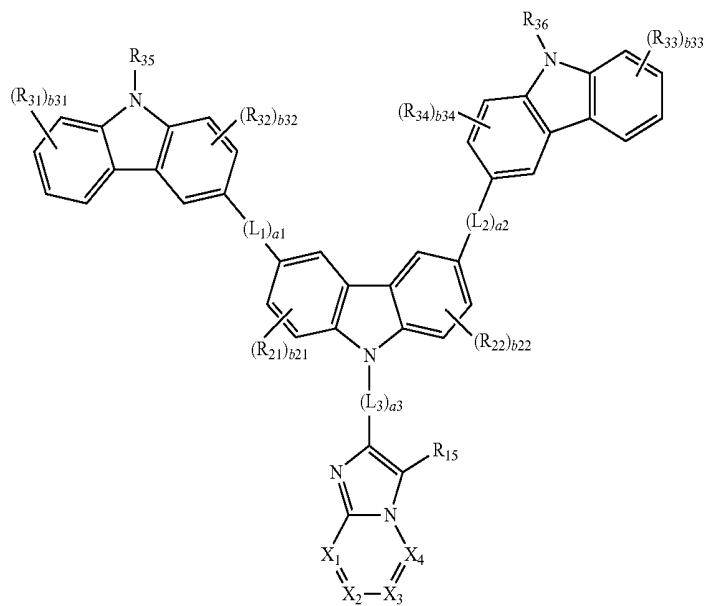
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<Formula 1A(5)>

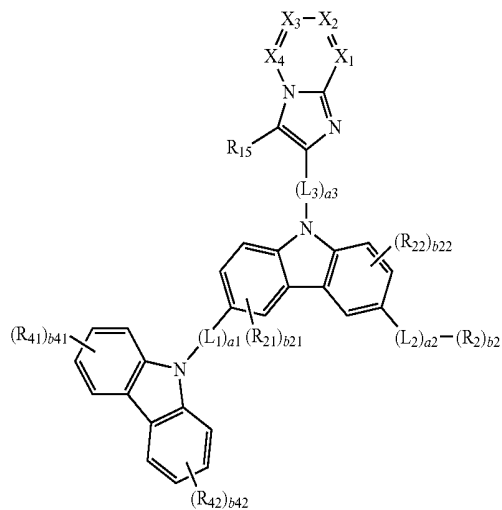


<Formula 1A(6)>

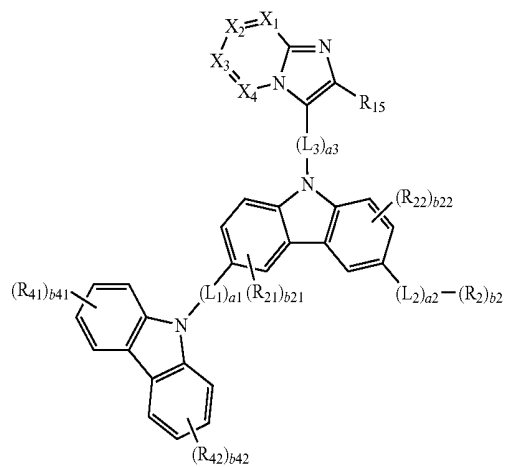


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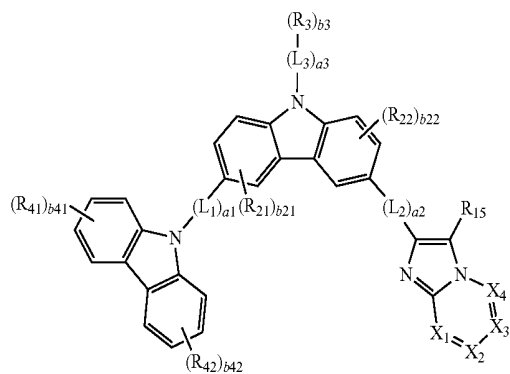
<Formula 1B(1)>



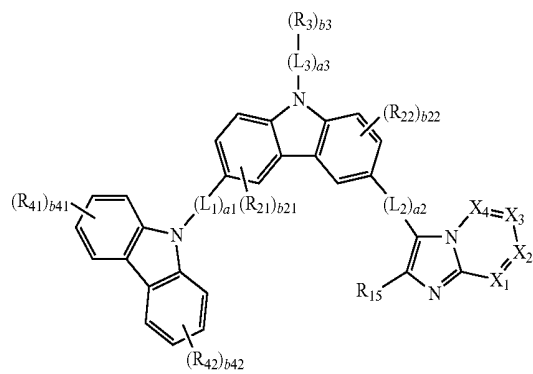
<Formula 1B(2)>



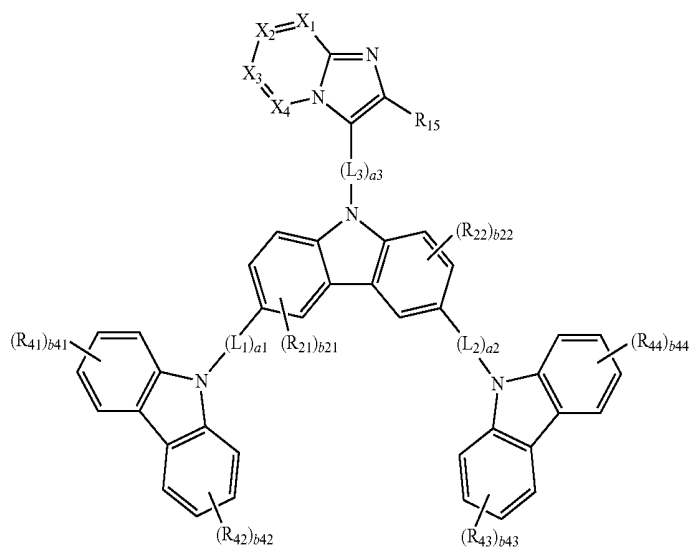
<Formula 1B(3)>



<Formula 1B(4)>

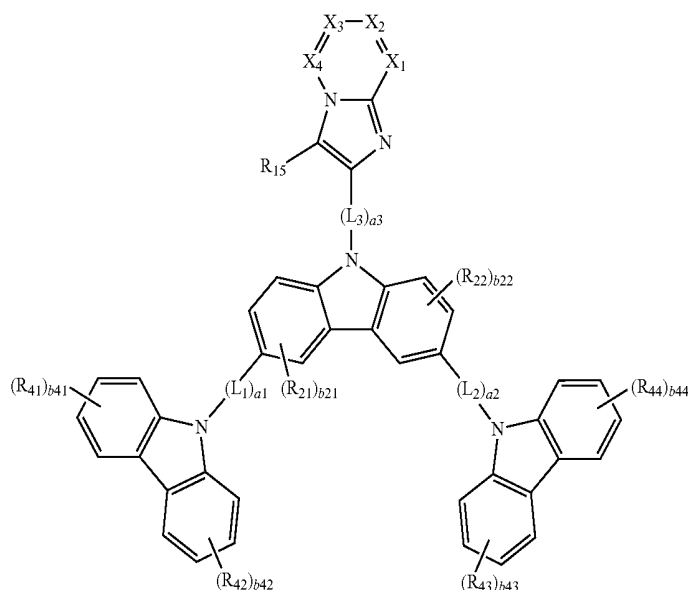


<Formula 1B(5)>



-continued

<Formula 1B(6)>



[0118] In Formulae 1A(1) to 1A(6) and Formulae 1B(1) to 1B(6),

[0119] X_1 to X_4 , R_{15} , R_{21} , R_{22} , b_{21} , b_{22} , L_1 to L_3 , a_1 to a_3 , b_2 and b_3 may be the same as defined herein;

[0120] R_2 and R_3 may be each independently selected from or include, e.g., a group represented by Formula 2, a hydrogen, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a substituted or unsubstituted C_1 - C_{60} alkyl group, a substituted or unsubstituted C_2 - C_{60} alkenyl group, a substituted or unsubstituted C_2 - C_{60} alkynyl group, a substituted or unsubstituted C_1 - C_{60} alkoxy group, a substituted or unsubstituted C_3 - C_{10} cycloalkyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkyl group, a substituted or unsubstituted C_3 - C_{10} cycloalkenyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenyl group, a substituted or unsubstituted C_6 - C_{60} aryl group, a substituted or unsubstituted C_1 - C_{60} heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group (wherein the "substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group" does not include the group represented by Formula 2 above), $-\text{Si}(Q_1)(Q_2)(Q_3)$, $-\text{B}(Q_4)(Q_5)$, and $-\text{N}(Q_6)(Q_7)$;

[0121] R_{31} to R_{34} and R_{41} to R_{44} are the same as R_{11} defined herein;

[0122] R_{35} and R_{36} may be each independently selected from or include, e.g., a substituted or unsubstituted C_1 - C_{60} alkyl group, a substituted or unsubstituted C_2 - C_{60} alkenyl group, a substituted or unsubstituted C_2 - C_{60} alkynyl group, a substituted or unsubstituted C_1 - C_{60} alkoxy group, a substituted or unsubstituted C_3 - C_{10} cycloalkyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkyl group, a substituted or unsubstituted C_3 - C_{10} cycloalkenyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenyl group, a substituted or unsubstituted C_6 - C_{60} aryl group, a substituted or unsubstituted C_1 - C_{60} heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, and a

substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group;

[0123] b_{31} , b_{33} , b_{41} to b_{44} may be each independently selected from 0, 1, 2, 3, and 4, and b_{32} and b_{34} may be each independently selected from 0, 1, 2, and 3.

[0124] For example, in Formulae 1A(1) to 1A(6) and Formulae 1B(1) to 1B(6),

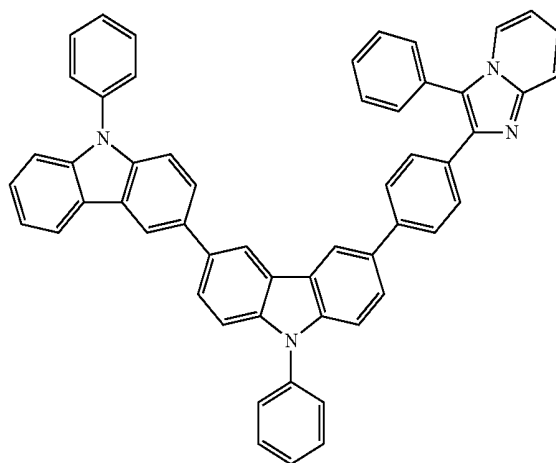
[0125] X_1 to X_4 may be CH; or

[0126] X_1 may be N, and X_2 to X_4 may be CH.

[0127] In some embodiments, a carbazole-based compound represented by Formula 1 may be represented by one selected from Formulae 1A(1), 1A(2), 1A(4), 1B(1) and 1B(2), but is not limited thereto.

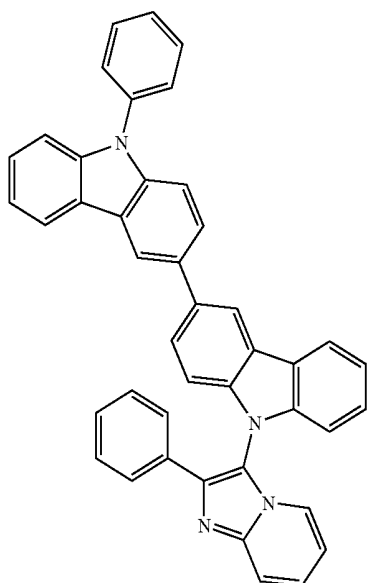
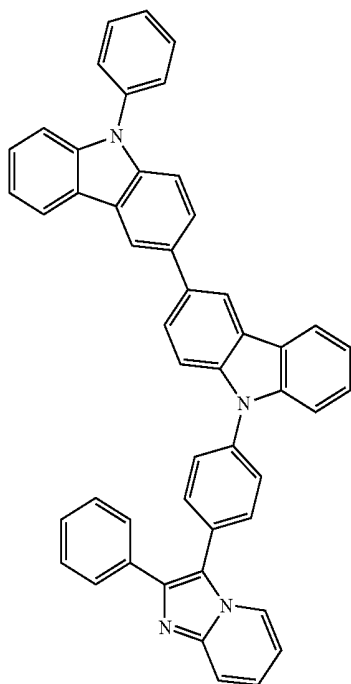
[0128] According to an embodiment, in Formulae 1A(1) to 1A(6) and Formulae 1B(1) to 1B(6), L_1 to L_3 may be each independently selected from a group represented by one of Formulae 4-1 to 4-24, and a_1 to a_3 may be each independently 0 or 1.

[0129] A carbazole-based compound represented by Formula 1 may be one selected from Compounds 1 to 20 below, but is not limited thereto.

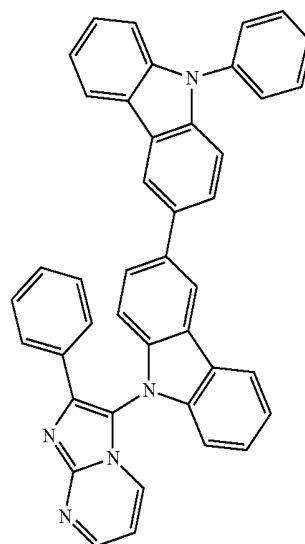
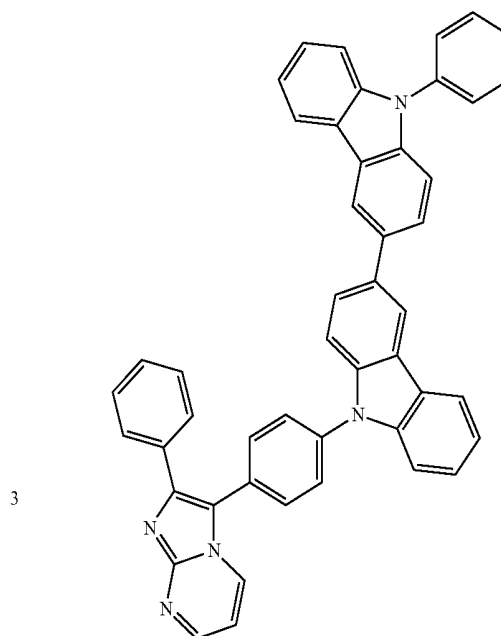
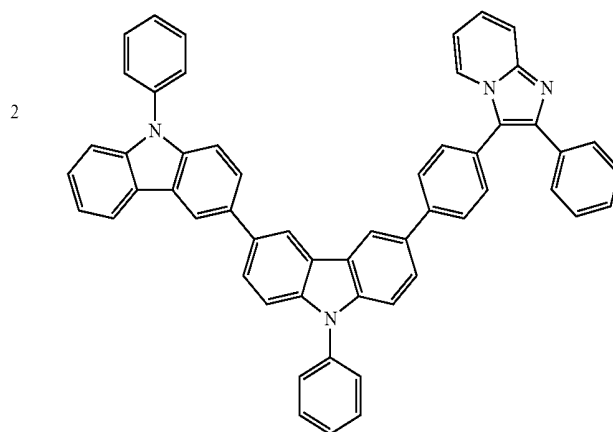


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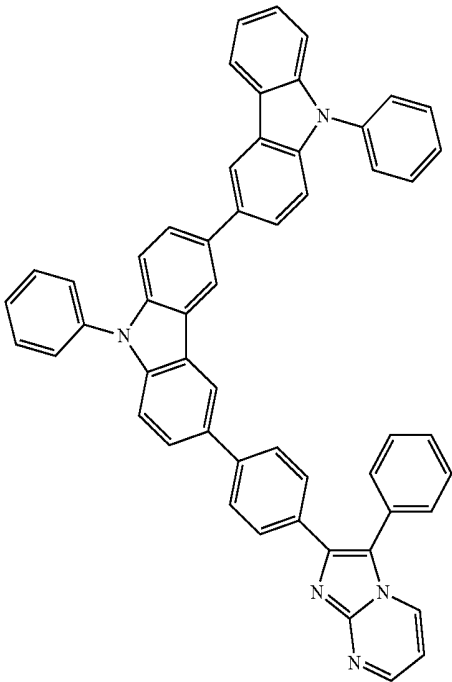


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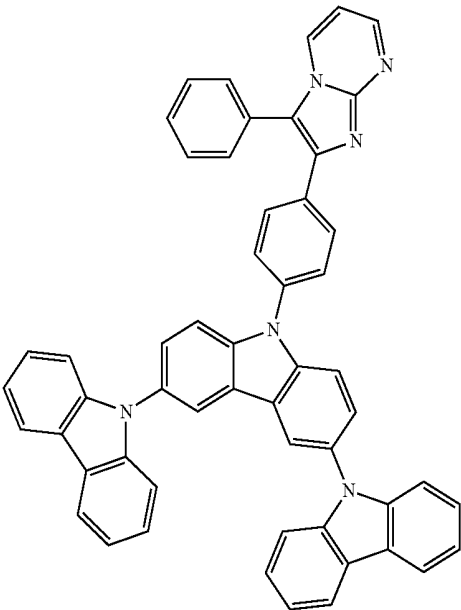
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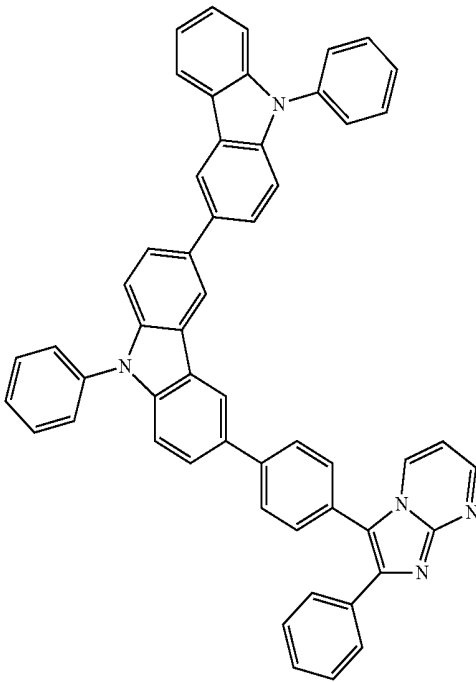


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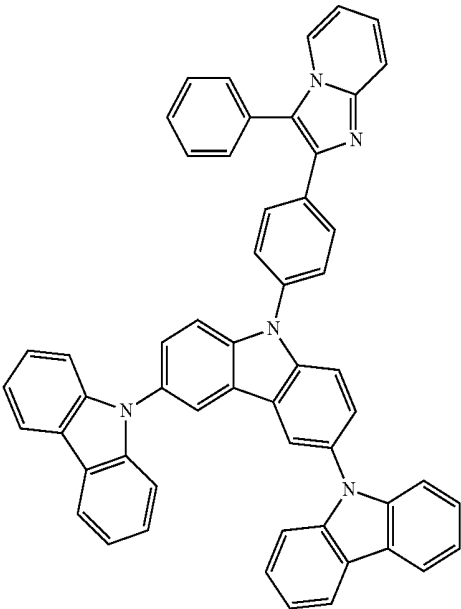
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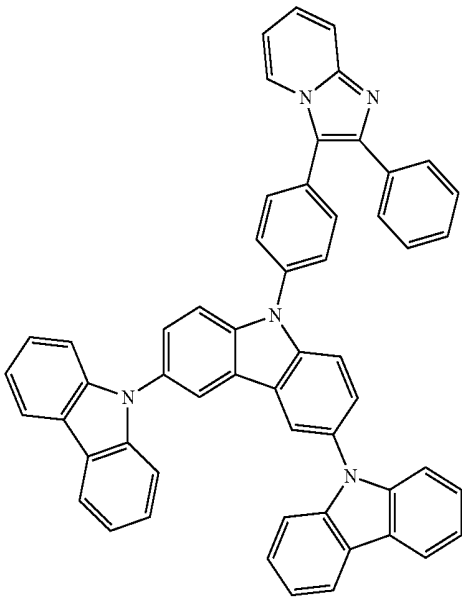


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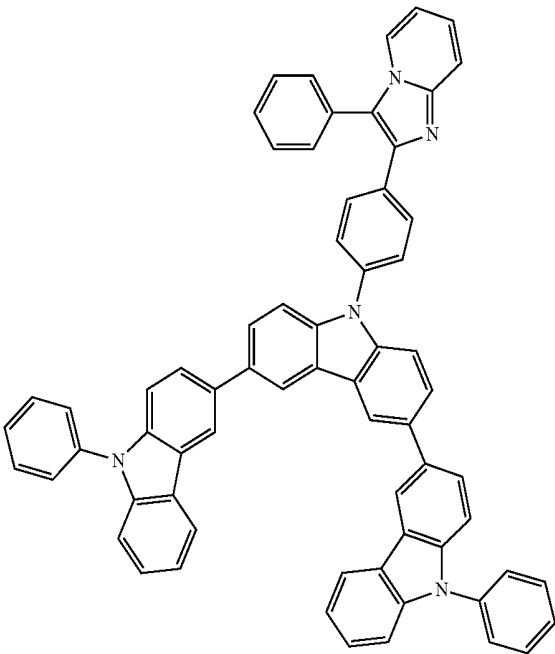
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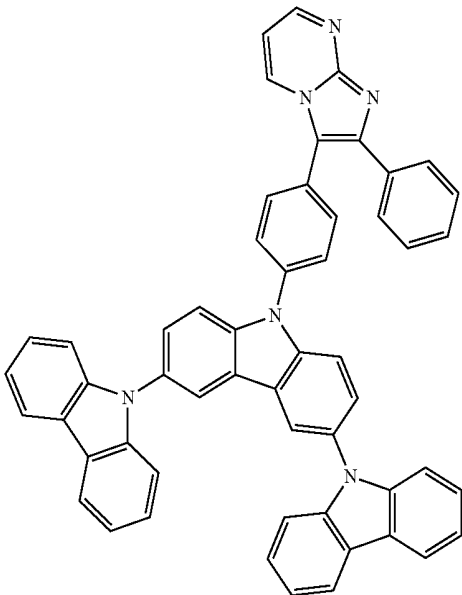


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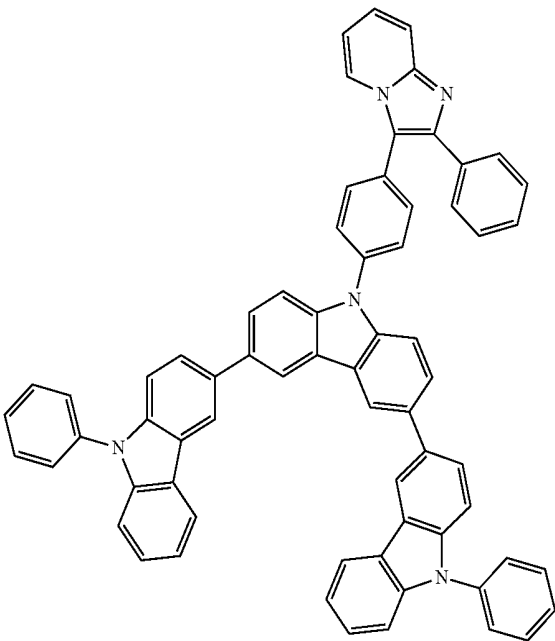
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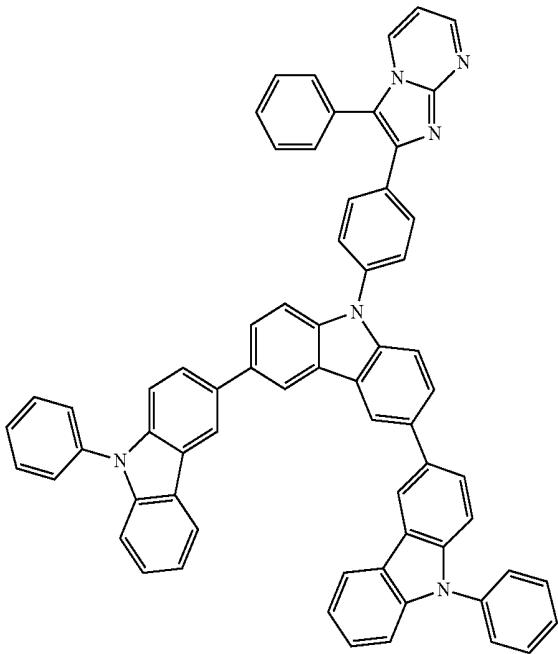


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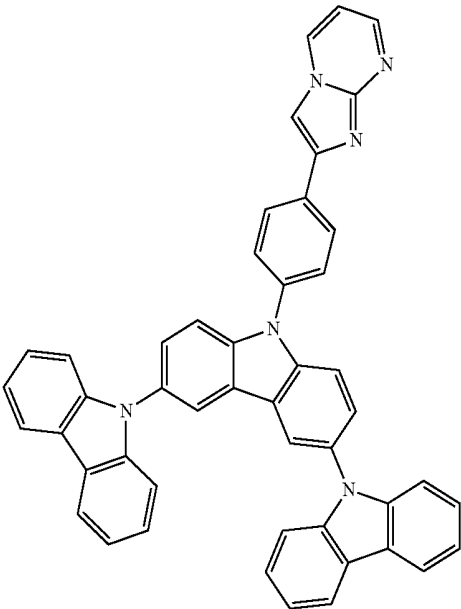
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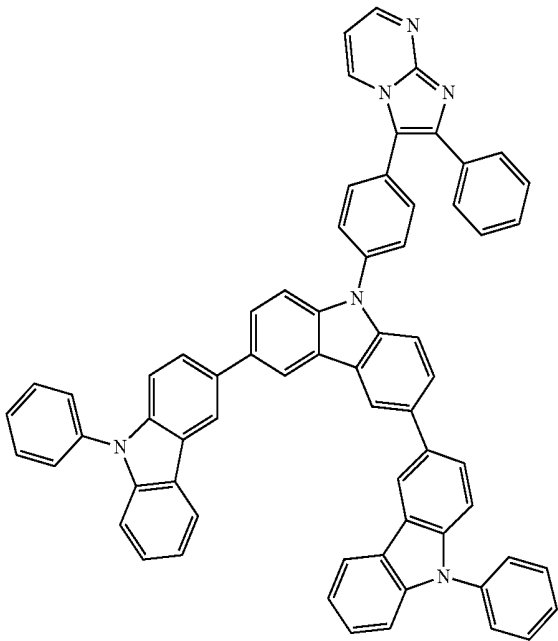


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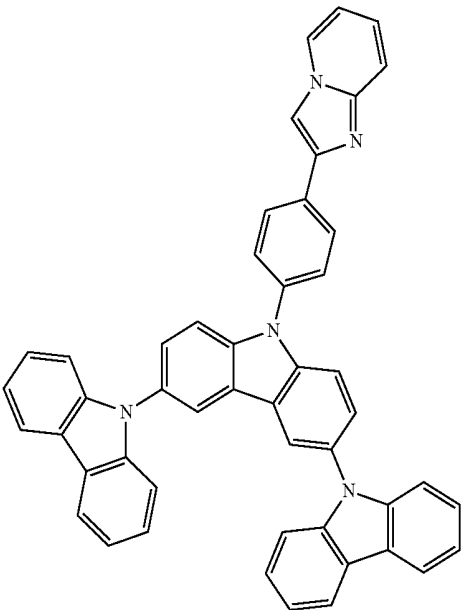
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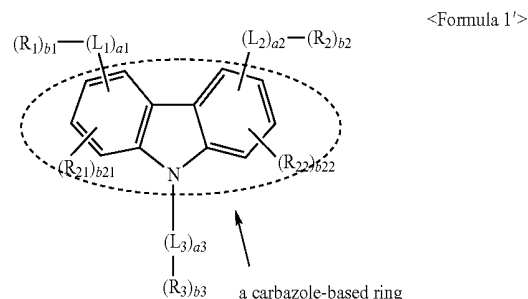
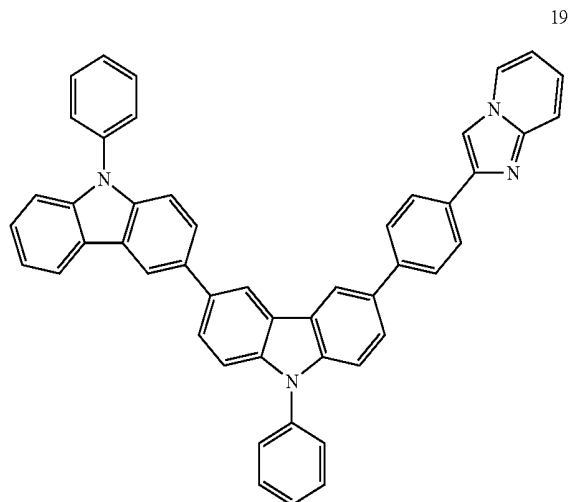
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[0131] Also, at least one of R_2 and R_3 in the carbazole-based compound represented by Formula 1 may be a group represented by Formula 2. b_2 and b_3 respectively indicate the number of R_2 and R_3 , and b_2 and b_3 may each independently be an integer of 1 to 3. For example, the carbazole-based ring represented by Formula 1' above is necessarily substituted with “the group represented by Formula 2”. The group represented by Formula 2, e.g., an imidazopyridinyl group or an imidazopyrimidinyl group, may have an excellent electron transporting characteristic, and thus the carbazole-based ring in Formula 1' substituted with one or more groups represented by Formula 2 may provide the compound represented by Formula 1 with an excellent electron injection characteristic. Accordingly, an organic light-emitting device including the groups represented by Formula 2 may have a low driving voltage.

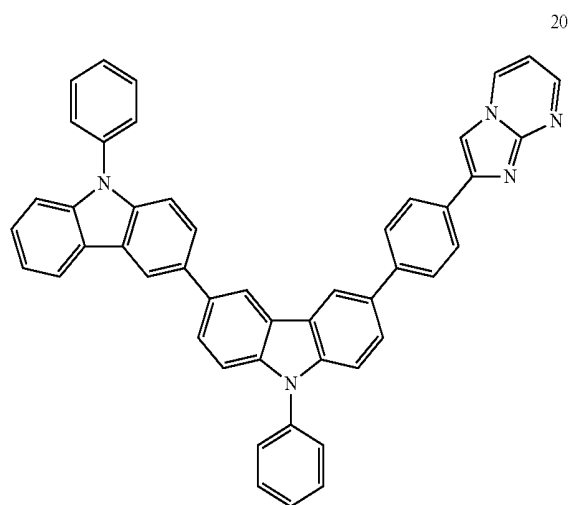
[0132] The carbazole-based compound represented by Formula 1 may be synthesized by using a suitable organic synthetic method. A synthesis method of the carbazole-based compound may be understood in view of the following embodiments.

[0133] At least one of the carbazole-based compounds represented by Formula 1 may be used or included between a pair of electrodes of an organic light-emitting device. In an implementation, the carbazole-based compound may be included in an electron transport region, e.g., an electron transport layer. Accordingly, an organic light-emitting device according to an embodiment may include: a first electrode; a second electrode facing the first electrode; and an organic layer that is disposed between the first electrode and the second electrode and includes an emission layer, wherein the organic layer includes at least one of the carbazole-based compounds described above.

[0134] As used herein, “(An organic layer) includes at least one of carbazole-based compounds” may be understood that “(An organic layer) may include one of carbazole-based compounds represented by Formula 1, or two or more different carbazole-based compounds represented by Formula 1”.

[0135] For example, the organic layer may include only Compound 1 as the carbazole-based compound. In this regard, Compound 1 may exist in an emission layer of the organic light-emitting device. In another embodiment, the organic layer may include, as the carbazole-based compound, Compound 1 and Compound 2. In this regard, Compound 1 and Compound 2 may exist in an identical layer (for example, Compound 1 and Compound 2 may both exist in an emission

[0130] The carbazole-based compound represented by Formula 1 may be or include a substituted or unsubstituted carbazolyl group. b indicates the number of R_1 , and b_1 is an integer of 1 to 3. For example, a carbazole-based ring represented by Formula 1' is necessarily substituted with “a substituted or unsubstituted carbazolyl group”. In this regard, the carbazole-based compound represented by Formula 1 may have a high T_i level suitable for a material of an organic light-emitting device, e.g., a green phosphorescent host. In an implementation, the carbazole-based compound represented by Formula 1 may have both an electron transporting substituent and a hole transporting substituent, and thus an organic light-emitting device including the carbazole-based compound may have a low driving voltage. (see Formula 1' below)



layer), or different layers (for example, Compound 1 may exist in an emission layer and Compound 2 may exist in an electron transport layer).

[0136] The organic layer may include, e.g., i) a hole transport region that is disposed between the first electrode and the emission layer and includes at least one selected from a hole injection layer, a hole transport layer, a buffer layer, and an electron blocking layer, and ii) an electron transport region that is disposed between the emission layer and the second electrode and includes at least one selected from a hole blocking layer, an electron transport layer, and an electron injection layer. In an implementation, the emission layer may include the carbazole-based compound represented by Formula 1.

[0137] The term "organic layer" used herein refers to a single layer and/or a plurality of layers disposed between the first electrode and the second electrode of an organic light-emitting device. A material included in the "organic layer" is not limited to an organic material.

[0138] FIG. 1 illustrates a schematic view of an organic light-emitting device 10 according to an embodiment. The organic light-emitting device 10 may include a first electrode 110, an organic layer 150, and a second electrode 190.

[0139] Hereinafter, the structure of an organic light-emitting device according to an embodiment and a method of manufacturing an organic light-emitting device according to an embodiment will be described in connection with FIG. 1.

[0140] In FIG. 1, a substrate may be additionally disposed under the first electrode 110 or above the second electrode 190. The substrate may be a glass substrate or transparent plastic substrate, each with excellent mechanical strength, thermal stability, transparency, surface smoothness, ease of handling, and water repellency.

[0141] The first electrode 110 may be formed by depositing or sputtering a material for forming the first electrode on the substrate. When the first electrode 110 is an anode, the material for the first electrode may be selected from materials with a high work function to make holes be easily injected. The first electrode 110 may be a reflective electrode or a transmissive electrode. The material for the first electrode may be a transparent and highly conductive material, and examples of such a material are indium tin oxide (ITO), indium zinc oxide (IZO), tin oxide (SnO_2), and zinc oxide (ZnO). When the first electrode 110 is a semi-transmissive electrode or a reflective electrode, as a material for forming the first electrode, at least one selected from magnesium (Mg), aluminum (Al), aluminum-lithium (Al—Li), calcium (Ca), magnesium-indium (Mg—In), magnesium-silver (Mg—Ag) may be used.

[0142] The first electrode 110 may have a single-layer structure, or a multi-layer structure including two or more layers. For example, the first electrode 110 may have a three-layer structure of ITO/Ag/ITO, but the structure of the first electrode 110 is not limited thereto.

[0143] The organic layer 150 is disposed on the first electrode 110. The organic layer 150 may include an emission layer.

[0144] The organic layer 150 may further include a hole transport region disposed between the first electrode and the emission layer, and an electron transport region disposed between the emission layer and the second electrode.

[0145] The hole transport region may include at least one selected from a hole injection layer (HIL), a hole transport

layer (HTL), a buffer layer, and an electron blocking layer (EBL), and the electron transport region may include at least one selected from a hole blocking layer (HBL), an electron transport layer (ETL), and an electron injection layer (EIL), but they are not limited thereto.

[0146] The hole transport region may have a single-layer structure formed of a single material, a single-layer structure formed of a plurality of different materials, or a multi-layer structure having a plurality of layers formed of a plurality of different materials.

[0147] For example, the hole transport region may have a single-layer structure formed of a plurality of different materials, or a structure of hole injection layer/hole transport layer, a structure of hole injection layer/hole transport layer/buffer layer, a structure of hole injection layer/buffer layer, a structure of hole transport layer/buffer layer, or a structure of hole injection layer/hole transport layer/electron blocking layer, wherein layers of each structure are sequentially stacked from the first electrode 110 in this stated order, but are not limited thereto.

[0148] When the hole transport region includes a hole injection layer, the hole injection layer may be formed on the first electrode 110 by using various methods, such as vacuum deposition, spin coating casting, a Langmuir-Blodgett (LB) method, ink-jet printing, laser-printing, or laser-induced thermal imaging.

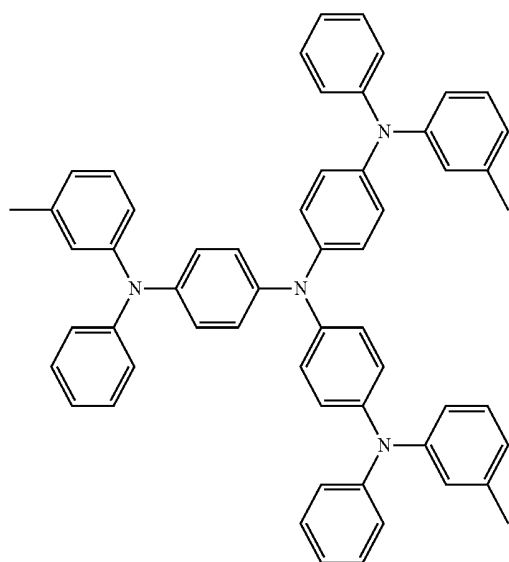
[0149] When a hole injection layer is formed by a vacuum deposition, for example, the vacuum deposition may be performed at a deposition temperature of about 100° C. to about 500° C., at a vacuum degree of about 10^{-8} torr to about 10^{-3} torr, and at a deposition rate of about 0.01 Å/sec to about 100 Å/sec in consideration of a compound for a hole injection layer to be deposited, and the structure of a hole injection layer to be formed.

[0150] When a hole injection layer is formed by a spin coating, the spin coating may be performed at a coating rate of about 2000 rpm to about 5000 rpm, and at a temperature of about 80° C. to 200° C. in consideration of a compound for a hole injection layer to be deposited, and the structure of a hole injection layer to be formed.

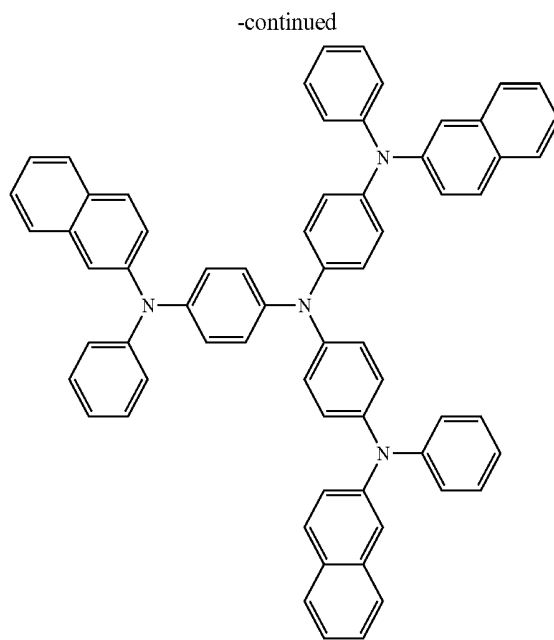
[0151] When the hole transport region includes a hole transport layer, the hole transport layer may be formed on the first electrode 110 or the hole injection layer by using various methods, such as vacuum deposition, spin coating, casting, a LB method, ink-jet printing, laser-printing, or laser-induced thermal imaging. When the hole transport layer is formed by vacuum deposition or spin coating, deposition and coating conditions for the hole transport layer may be determined by referring to the deposition and coating conditions for the hole injection layer.

[0152] The hole transport region may include at least one selected from m-MTDATA, TDATA, 2-TNATA, NPB, β -NPB, TPD, Spiro-TPD, Spiro-NPB, α -NPB, TAPC, HMTDP, 4,4',4''-tris(N-carbazolyl)triphenylamine (TCTA), polyaniline/dodecylbenzenesulfonic acid (Pani/DBSA), poly(3,4-ethylenedioxythiophene)/poly(4-styrenesulfonate) (PEDOT/PSS), polyaniline/camphor sulfonic acid (PANI/CSA),

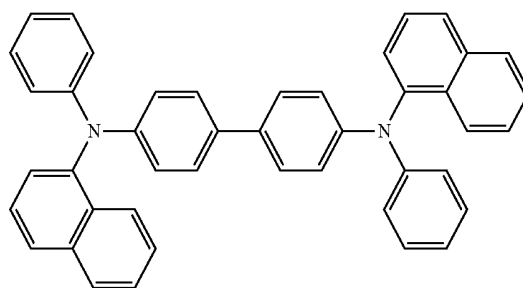
(polyaniline)/poly(4-styrenesulfonate) (Pani/PSS), a compound represented by Formula 201 below, and a compound represented by Formula 202 below:



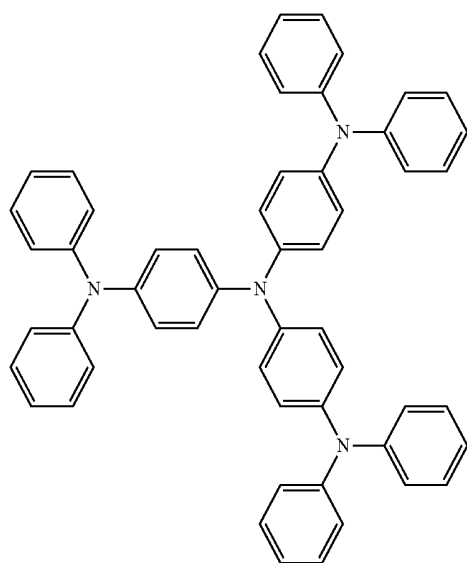
m-MTDATA



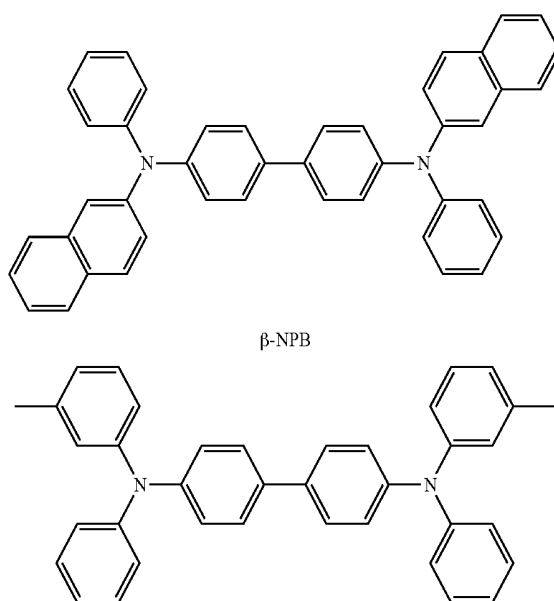
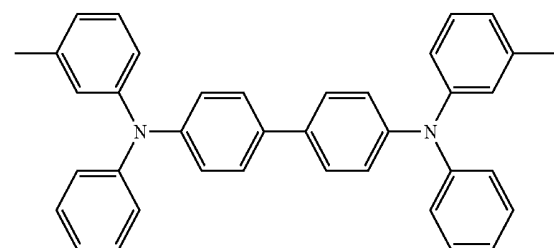
2-TNATA



NPB

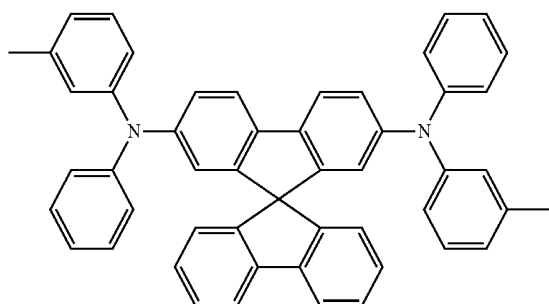


TDATA

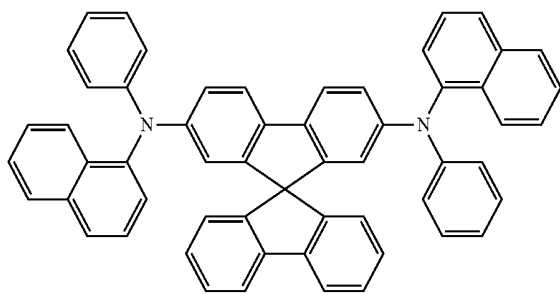
 β -NPB

TPD

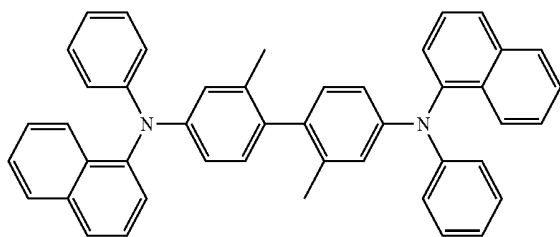
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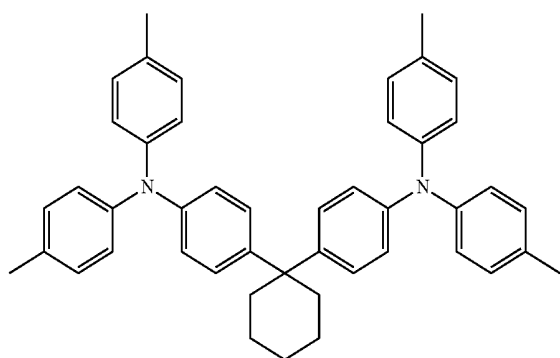
Spiro-TPD



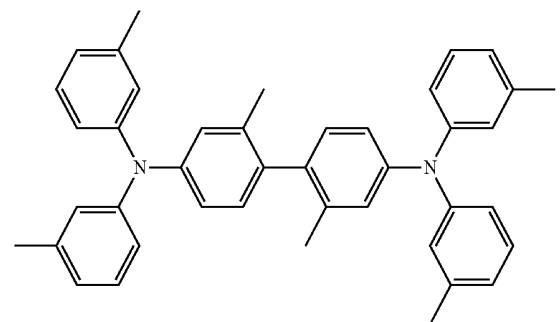
Spiro-NPB



methylated NPB

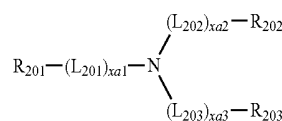


TAPC

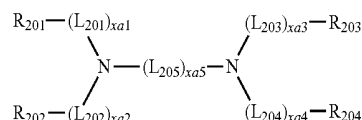


HMTPD

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<Formula 201>



<Formula 202>

[0153] wherein in Formulae 201 and 202,

[0154] L_{201} to L_{205} may be the same as explained in connection with L_1 ;[0155] xa1 to xa4 may be each independently selected from 0, 1, 2, and 3;[0156] xa5 may be selected from 1, 2, 3, 4, and 5; and[0157] R_{201} to R_{204} may be selected from a substituted or unsubstituted C_3 - C_{10} cycloalkyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkyl group, a substituted or unsubstituted C_3 - C_{10} cycloalkenyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenyl group, a substituted or unsubstituted C_6 - C_{60} aryl group, a substituted or unsubstituted C_6 - C_{60} aryloxy group, a substituted or unsubstituted C_6 - C_{60} arylthio group, a substituted or unsubstituted C_1 - C_{60} heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, and a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group.

[0158] For example, in Formulae 201 and 202,

[0159] L_{201} to L_{205} may be each independently selected from

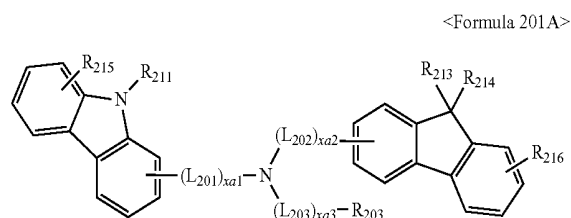
[0160] a phenylene group, a naphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzofluorene group, a dibenzofluorene group, a phenanthrenylene group, an anthracenylenylene group, a pyrenylene group, a chrysenylene group, a pyridinylenylene group, a pyrazinylenylene group, a pyrimidinylenylene group, a pyridazinylenylene group, a quinolinylenylene group, an isoquinolinylenylene group, a quinoxalinylenylene group, a quinazolinylenylene group, a carbazolinylenylene group and a triazinylenylene group; and

[0161] a phenylene group, a naphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzofluorene group, a dibenzofluorene group, a phenanthrenylene group, an anthracenylenylene group, a pyrenylene group, a chrysenylene group, a pyridinylenylene group, a pyrazinylenylene group, a pyrimidinylenylene group, a pyridazinylenylene group, a quinolinylenylene group, an isoquinolinylenylene group, a quinoxalinylenylene group, a quinazolinylenylene group, a carbazolinylenylene group, and a triazinylenylene group, each substituted with at least one selected from a deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazoliny group, a carbazolyl group, and a triazinyl group;[0162] xa1 to xa4 may be each independently 0, 1, or 2;[0163] xa5 may be 1, 2, or 3; and[0164] R_{201} to R_{204} are each independently selected from

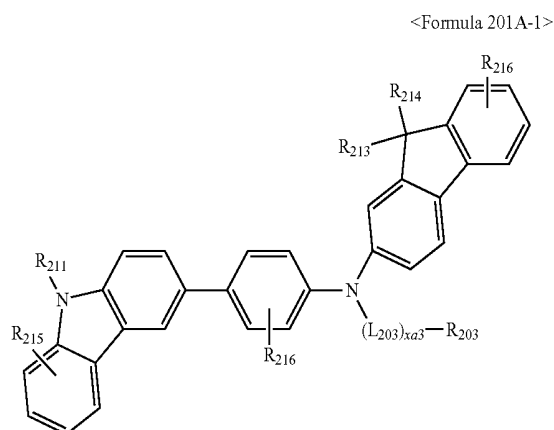
[0165] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group and a triazinyl group; and

[0166] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, and a triazinyl group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, a naphthyl group, an azulenyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, and a triazinyl group, but they are not limited thereto.

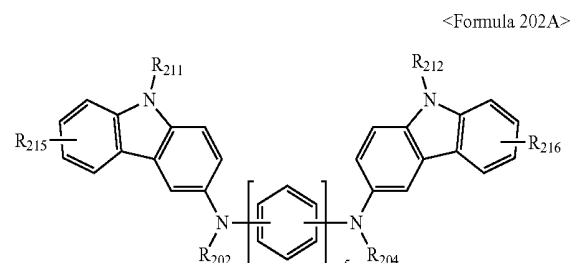
[0167] The compound represented by Formula 201 may be represented by Formula 201A:



[0168] For example, the compound represented by Formula 201 may be represented by Formula 201A-1 below, but is not limited thereto:



[0169] For example, the compound represented by Formula 202 may be represented by Formula 202A below, but is not limited thereto:



[0170] L₂₀₁ to L₂₀₃, xa1 to xa3, xa5, and R₂₀₂ to R₂₀₄ in Formulae 201A, 201A-1, and 202A are already described above, R₂₁₁ is the same as explained in connection with R₂₀₃, and R₂₁₃ to R₂₁₆ may be each independently selected from a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid group or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monvalent non-aromatic condensed polycyclic group, and a monvalent non-aromatic condensed heteropolycyclic group.

[0171] For example, in Formulae 201A, 201A-1, and 202A,

[0172] L₂₀₁ to L₂₀₃ may be each independently selected from

[0173] a phenylene group, a naphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzofluorenylene group, a dibenzofluorenylene group, a phenanthrenylene group, an anthracenylenylene group, a pyrenylene group, a chrysenylene group, a pyridinylenylene group, a pyrazinylenylene group, a pyrimidinylenylene group, a pyridazinylene group, a quinolinylene group, an isoquinolinylene group, a quinoxalinylenylene group, a quinazolinylene group, a carbazolylenylene group and a triazinylene group; and

[0174] a phenylene group, a naphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzofluorenylene group, a dibenzofluorenylene group, a phenanthrenylene group, an anthracenylenylene group, a pyrenylene group, a chrysenylene group, a pyridinylenylene group, a pyrazinylenylene group, a pyrimidinylenylene group, a pyridazinylene group, a quinolinylene group, an isoquinolinylene group, a quinoxalinylenylene group, a quinazolinylene group, a carbazolylenylene group, and a triazinylene group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzo-fluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl

group, an isoquinolinyl group, a quinoxaliny group, a quinazolinyl group, a carbazolyl group, and a triazinyl group; and

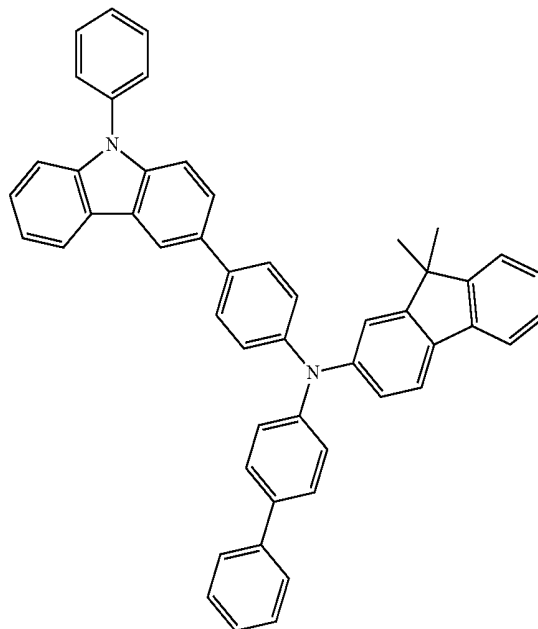
[0189] xa5 may be 1 or 2.

[0190] R₂₁₃ and R₂₁₄ in Formulae 201A, and 201A-1 may bind to each other to form a saturated or unsaturated ring.

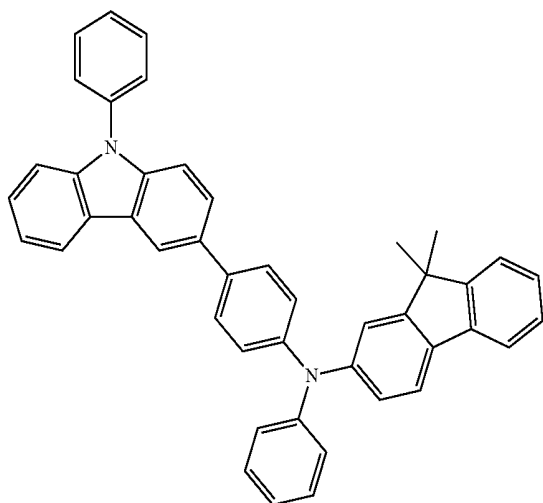
[0191] The compound represented by Formula 201, and the compound represented by Formula 202 may each include compounds HT1 to HT20 illustrated below, but are not limited thereto.

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HT3

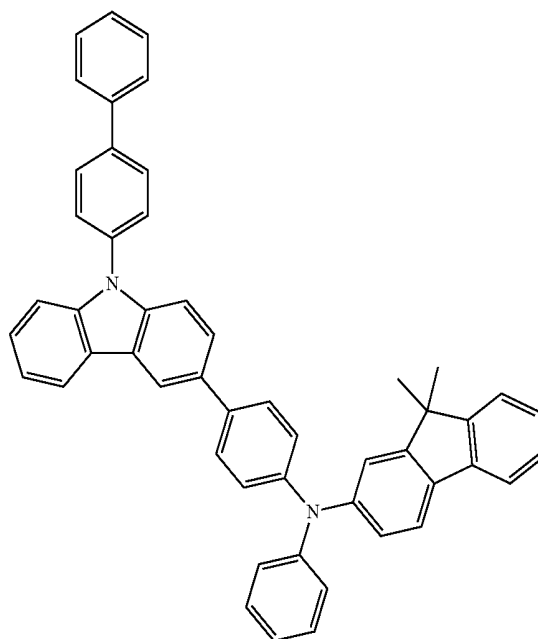
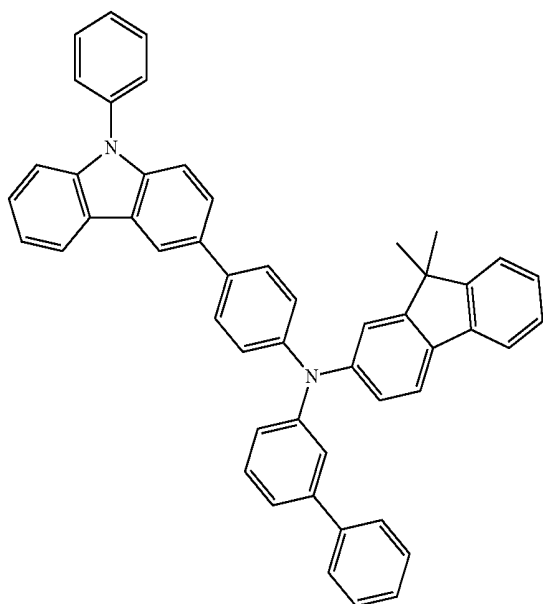


HT1



HT4

HT2

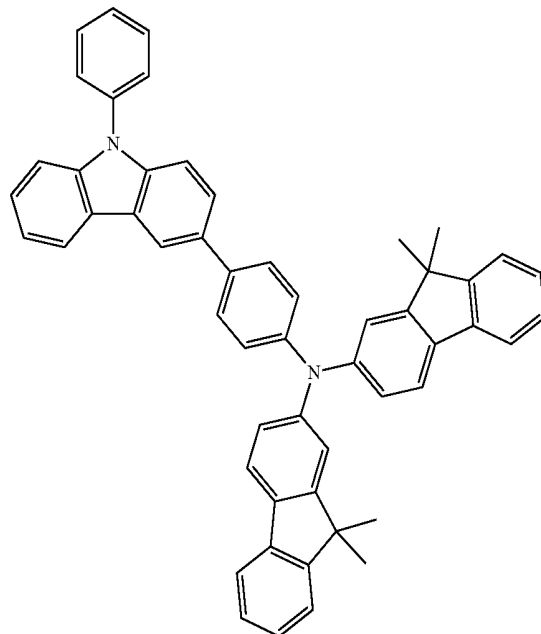
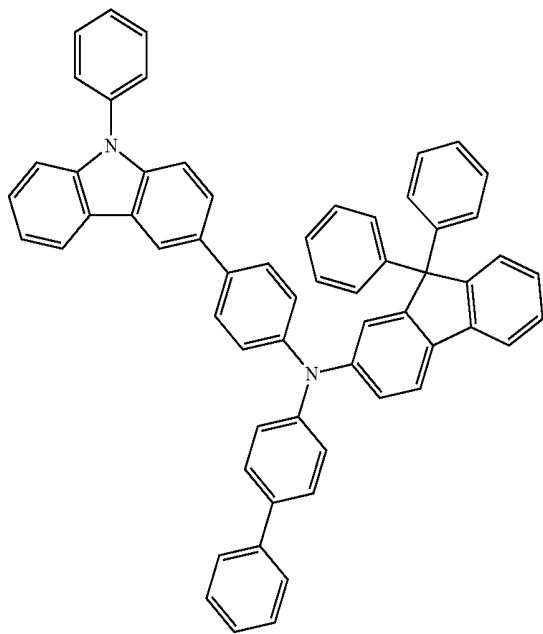


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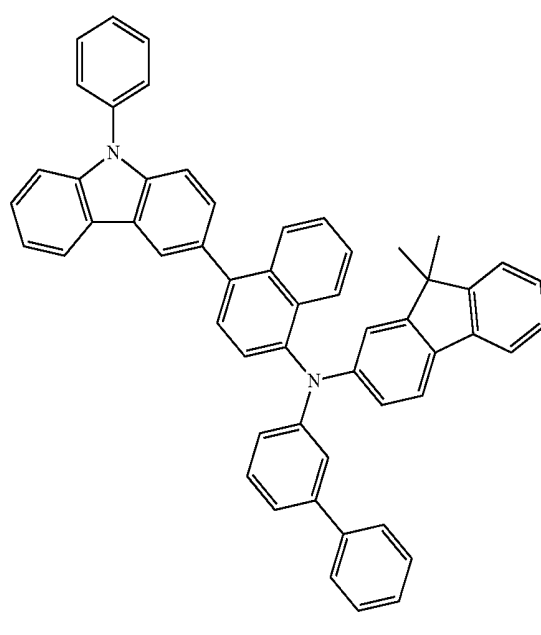
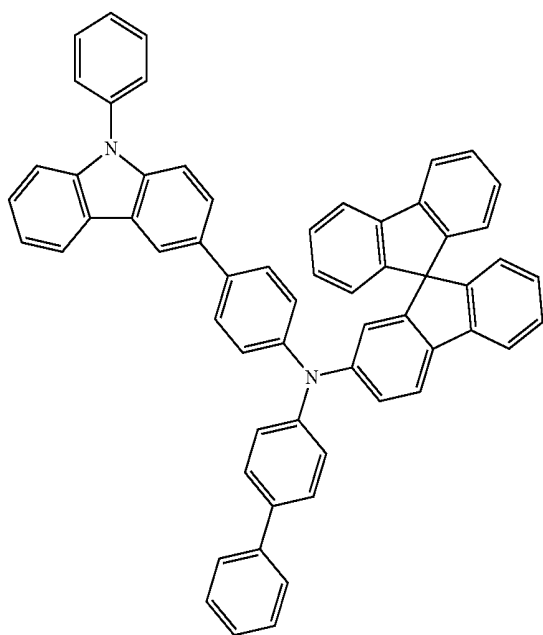
HT5

HT7



HT6

HT8

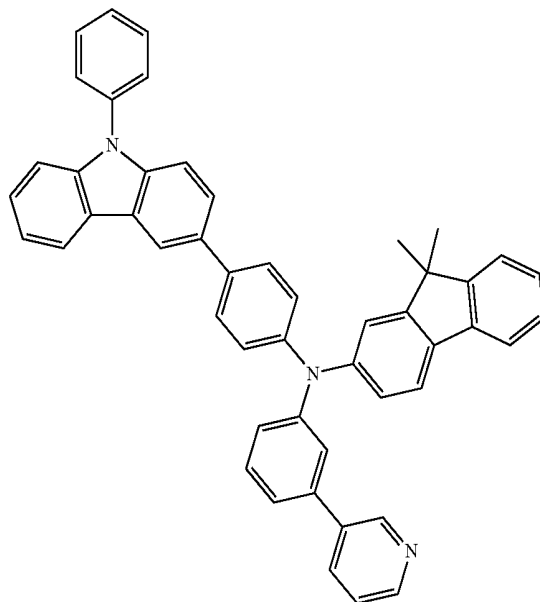
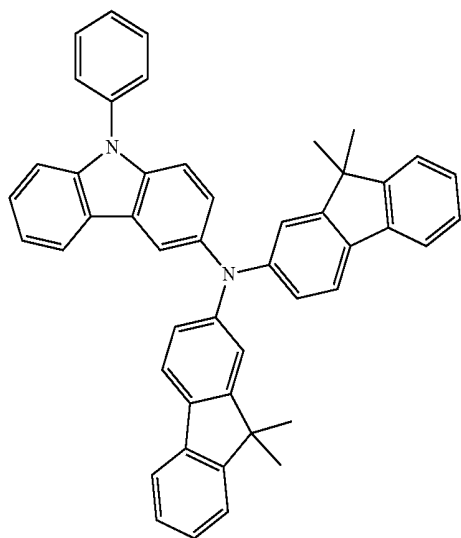


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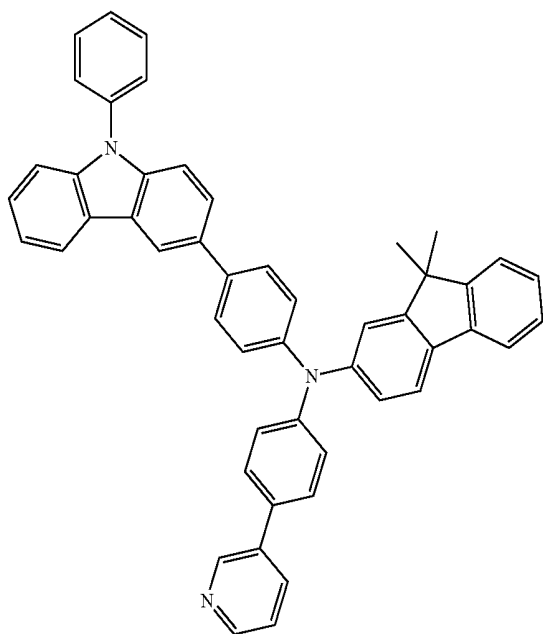
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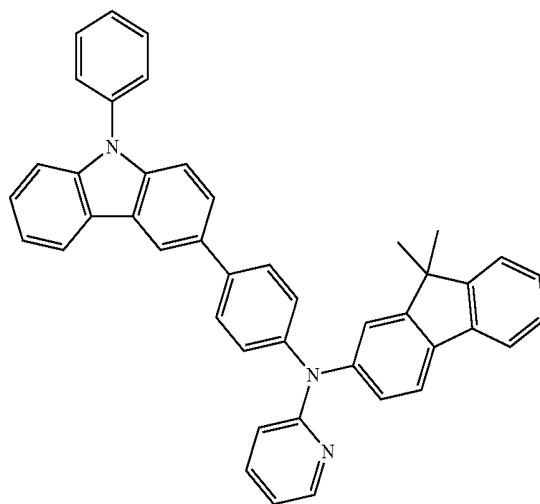
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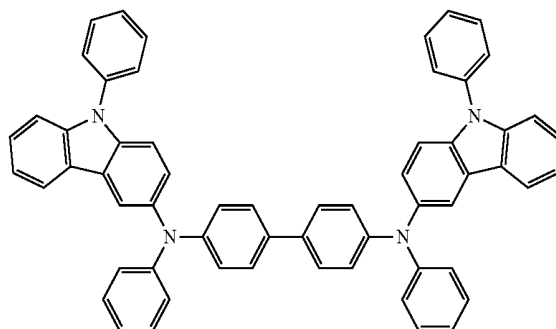
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HT12

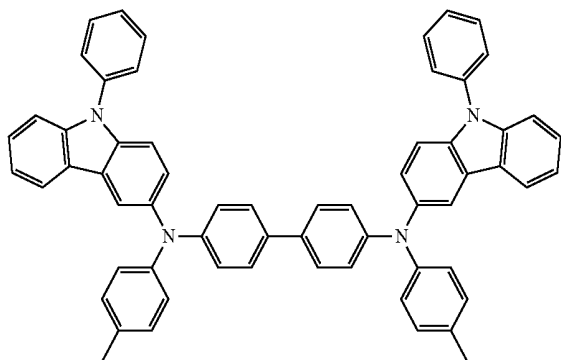


HT13



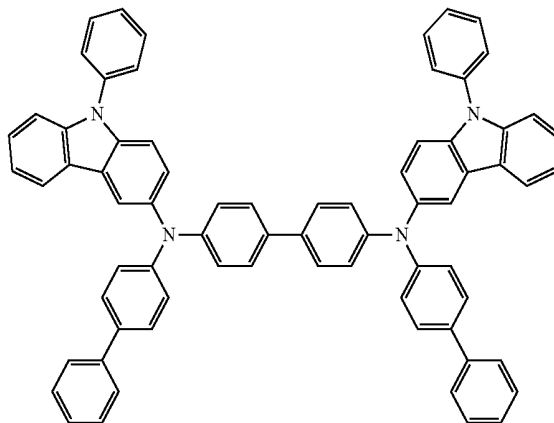
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HT14



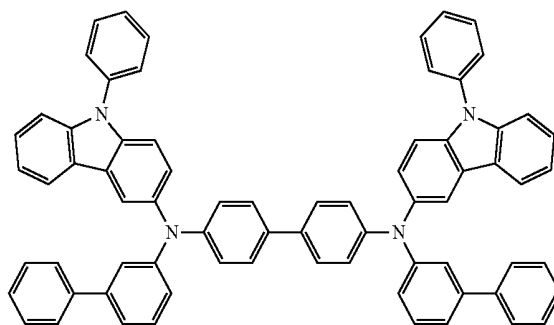
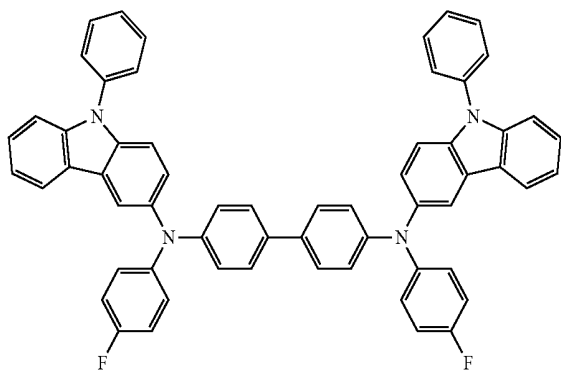
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HT18



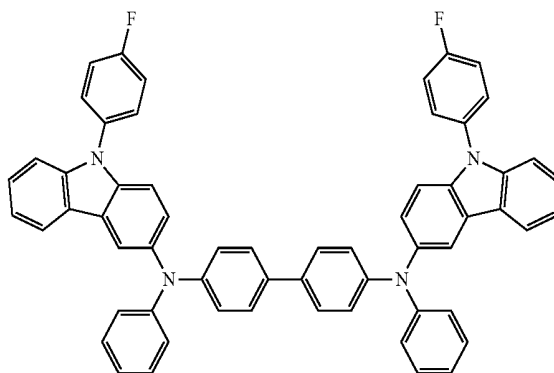
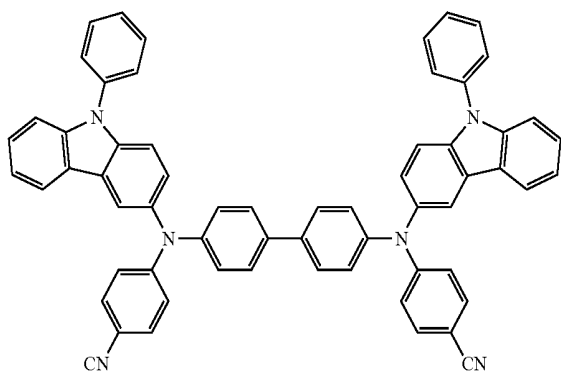
HT19

HT15

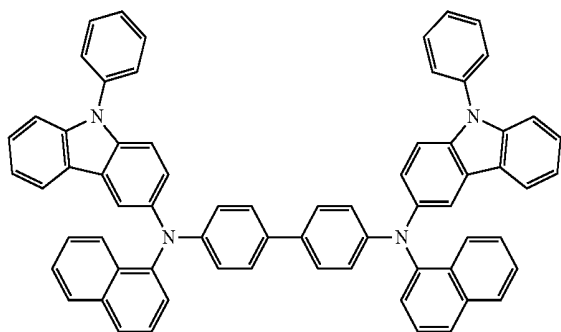


HT20

HT16



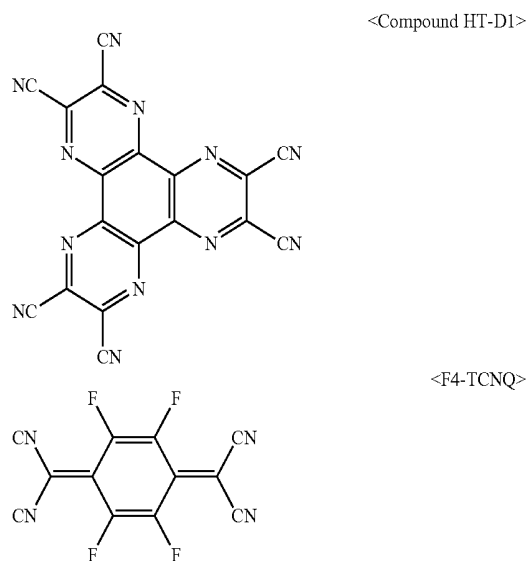
HT17



[0192] A thickness of the hole transport region may be in a range of about 100 Å to about 10000 Å, for example, about 100 Å to about 1000 Å. When the hole transport region includes a hole injection layer and a hole transport layer, a thickness of the hole injection layer may be in a range of about 100 Å to about 10000 Å, and for example, about 100 Å to about 1000 Å, and a thickness of the hole transport layer may be in a range of about 50 Å to about 2000 Å, for example, about 100 Å to about 1500 Å. When the thicknesses of the hole transport region, the hole injection layer, and the hole transport layer are within these ranges, satisfactory hole transporting characteristics may be obtained without a substantial increase in driving voltage.

[0193] The hole transport region may further include, in addition to these materials, a charge-generation material for the improvement of conductive properties. The charge-generation material may be homogeneously or non-homogeneously dispersed in the hole transport region.

[0194] The charge-generation material may be, e.g., a p-dopant. The p-dopant may be one selected from a quinone derivative, a metal oxide, and a cyano group-containing compound, but is not limited thereto. For example, non-limiting examples of the p-dopant are a quinone derivative, such as tetracyanoquinonodimethane (TCNQ) or 2,3,5,6-tetrafluoro-tetracyano-1,4-benzoquinonodimethane (F4-TCNQ); a metal oxide, such as a tungsten oxide or a molybdenum oxide, and Compound HT-D1 illustrated below, but are not limited thereto.



[0195] The hole transport region may further include, in addition to the hole injection layer and the hole transport layer, at least one selected from a buffer layer and an electron blocking layer. Since the buffer layer may compensate for an optical resonance distance according to a wavelength of light emitted from the emission layer, luminescent efficiency of a formed organic light-emitting device may be improved. For use as a material included in the buffer layer, materials that are included in the hole transport region may be used. The electron blocking layer prevents injection of electrons from the electron transport region.

[0196] An emission layer may be formed on the first electrode 110 or the hole transport region by using various methods, e.g., vacuum deposition, spin coating, casting, a LB method, ink-jet printing, laser-printing, or laser-induced thermal imaging. When the emission layer is formed by vacuum deposition or spin coating, deposition and coating conditions for the emission layer may be determined by referring to the deposition and coating conditions for the hole injection layer.

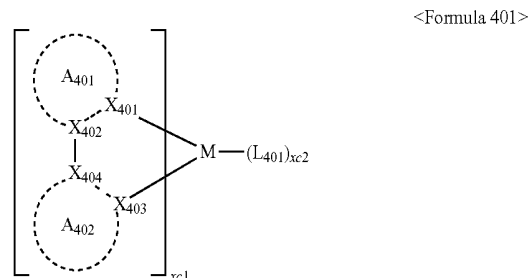
[0197] When the organic light-emitting device 10 is a full color organic light-emitting device, the emission layer may be patterned into a red emission layer, a green emission layer, or a blue emission layer for each sub pixel within the organic light-emitting device 10. In an implementation, the emission layer may have a stacked structure of a red emission layer, a green emission layer, and a blue emission layer, or may include a red-light emission material, a green-light emission material, and a blue-light emission material, which are mixed with each other in a single layer, to emit white light.

[0198] The emission layer may include the carbazole-based compound represented by Formula 1.

[0199] The emission layer may include a host and a dopant. The host may include the carbazole-based compound represented by Formula 1.

[0200] The dopant may be at least one selected from a fluorescent dopant and a phosphorescent dopant.

[0201] The phosphorescent dopant may include an organo-metallic complex represented by Formula 401 below.



[0202] In Formula 401,

[0203] M may be selected from iridium (Ir), platinum (Pt), osmium (Os), titanium (Ti), zirconium (Zr), hafnium (Hf), europium (Eu), terbium (Tb), and thulium (Tm);

[0204] X₄₀₁ to X₄₀₄ may be each independently nitrogen or carbon;

[0205] A₄₀₁ and A₄₀₂ rings may be each independently selected from or include, e.g., a substituted or unsubstituted benzene, a substituted or unsubstituted naphthalene, a substituted or unsubstituted fluorene, a substituted or unsubstituted spiro-fluorene, a substituted or unsubstituted indene, a substituted or unsubstituted pyrrole, a substituted or unsubstituted thiophene, a substituted or unsubstituted furan, a substituted or unsubstituted imidazole, a substituted or unsubstituted pyrazole, a substituted or unsubstituted thiazole, a substituted or unsubstituted isothiazole, a substituted or unsubstituted oxazole, a substituted or unsubstituted isoxazole, a substituted or unsubstituted pyridine, a substituted or unsubstituted pyrazine, a substituted or unsubstituted pyrimidine, a substituted or unsubstituted pyridazine, a substituted or unsubstituted quinoline, a substituted or unsubstituted isoquinoline, a substituted or unsubstituted benzoquinoline, a substituted or unsubstituted quinoxaline, a substituted or unsubstituted quinazoline, a substituted or unsubstituted carbazole, a substituted or unsubstituted benzimidazole, a substituted or unsubstituted benzofuran, a substituted or unsubstituted benzothiophene, a substituted or unsubstituted isobenzothiophene, a substituted or unsubstituted benzoxazole, a substituted or unsubstituted isobenzoxazole, a substituted or unsubstituted triazole, a substituted or unsubstituted oxadiazole, a substituted or unsubstituted triazine, a substituted or unsubstituted dibenzofuran, and a substituted or unsubstituted dibenzothiophene; and

[0206] at least one of substituents of the substituted benzene, substituted naphthalene, substituted fluorene, substituted spiro-fluorene, substituted indene, substituted pyrrole, substituted thiophene, substituted furan, substituted imidazole, substituted pyrazole, substituted thiazole, substituted isothiazole, substituted oxazole, substituted isoxazole, substituted pyridine, substituted pyrazine, substituted pyrimidine, substituted pyridazine, substituted quinoline, substituted isoquinoline, substituted benzoquinoline, substituted quinoxaline, substituted quinazoline, substituted carbazole, substituted benzimidazole, substituted benzofuran, substi-

tuted benzothiophene, substituted isobenzothiophene, substituted benzoxazole, substituted isobenzoxazole, substituted triazole, substituted oxadiazole, substituted triazine, substituted dibenzofuran, and substituted dibenzothiophene may be selected from

[0207] a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group;

[0208] a C₁-C₆₀ alkyl group, a C₁-C₆₀ alkenyl group, a C₁-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid group or a salt thereof, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, —N(Q₄₀₁)(Q₄₀₂), —Si(Q₄₀₃)(Q₄₀₄)(Q₄₀₅), and —B(Q₄₀₆)(Q₄₀₇);

[0209] a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic heterocondensed polycyclic group;

[0210] a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic heterocondensed polycyclic group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid group or a salt thereof, a C₁-C₆₀ alkyl group, a C₁-C₆₀ alkenyl group, a C₁-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic heterocondensed polycyclic group, —N(Q₄₁₁)(Q₄₁₂), —Si(Q₄₁₃)(Q₄₁₄)(Q₄₁₅), and —B(Q₄₁₆)(Q₄₁₇); and

[0211] —N(Q₄₂₁)(Q₄₂₂), —Si(Q₄₂₃)(Q₄₂₄)(Q₄₂₅), and —B(Q₄₂₆)(Q₄₂₇); and

[0212] L₄₀₁ may be an organic ligand;

[0213] xc1 may be 1, 2, or 3; and

[0214] xc2 may be 0, 1, 2, or 3.

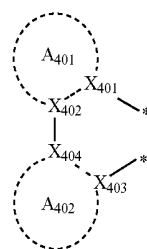
[0215] L₄₀₁ may be a monovalent, divalent, or trivalent organic ligand. For example, L₄₀₁ may be selected from a halogen ligand (for example, Cl or F), a diketone ligand (for example, acetylacetonate, 1,3-diphenyl-1,3-propanedionate,

2,2,6,6-tetramethyl-3,5-heptanedionate, or hexafluoroacetate), a carboxylic acid ligand (for example, picolinate, dimethyl-3-pyrazolecarboxylate, or benzoate), a carbon monooxide ligand, an isonitrile ligand, a cyano ligand, and a phosphorous ligand (for example, phosphine, and phosphite), but is not limited thereto.

[0216] When A₄₀₁ in Formula 401 has two or more substituents, the substituents of A₄₀₁ may bind to each other to form a saturated or unsaturated ring.

[0217] When A₄₀₂ in Formula 402 has two or more substituents, the substituents of A₄₀₂ may bind to each other to form a saturated or unsaturated ring.

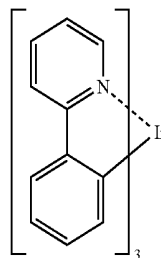
[0218] When xc1 in Formula 401 is two or more, a plurality of ligands



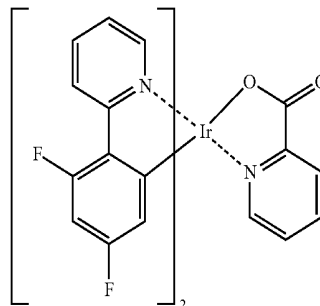
in Formula 401 may be identical or different. When xc1 in Formula 401 is two or more, A₄₀₁ and A₄₀₂ may be respectively directly connected to A₄₀₁ and A₄₀₂ of other neighboring ligands with or without a linker (for example, a C₁-C₅alkylene, or —N(R')— (wherein R' may be a C₁-C₁₀ alkyl group or a C₆-C₂₀ aryl group) or —C(=O)—) therebetween.

[0219] The phosphorescent dopant may include at least one selected from Compounds PD1 to PD74 below, but is not limited thereto.

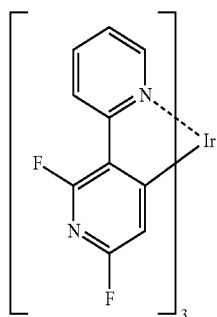
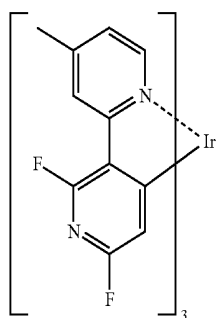
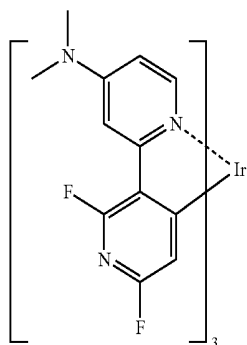
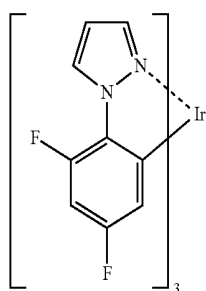
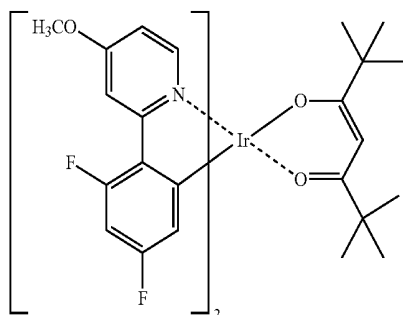
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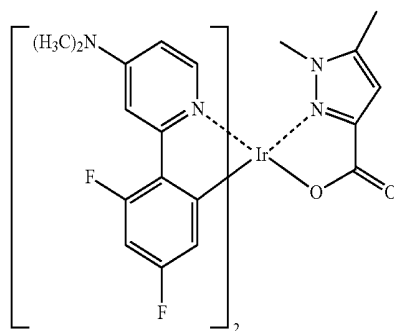
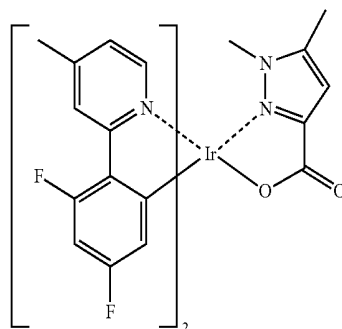
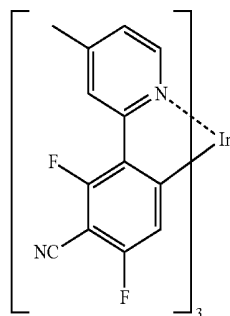
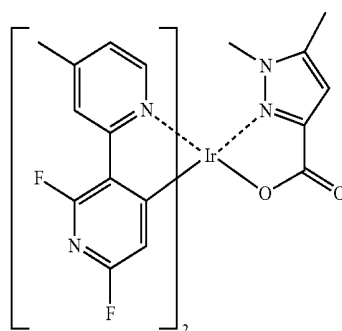
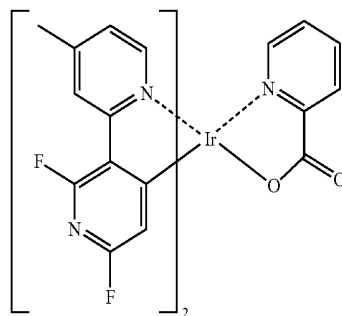
PD2



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PD3

PD8

PD4

PD9

PD5

PD10

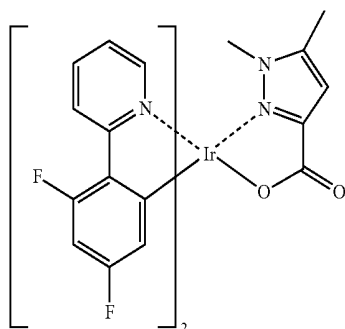
PD6

PD11

PD7

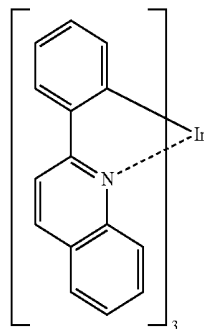
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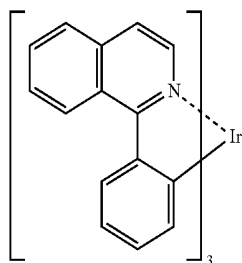


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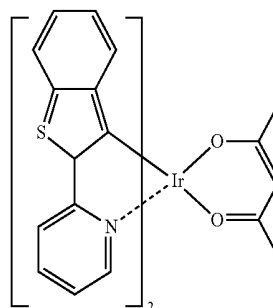
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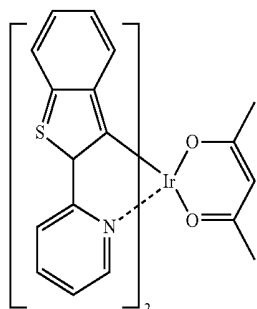
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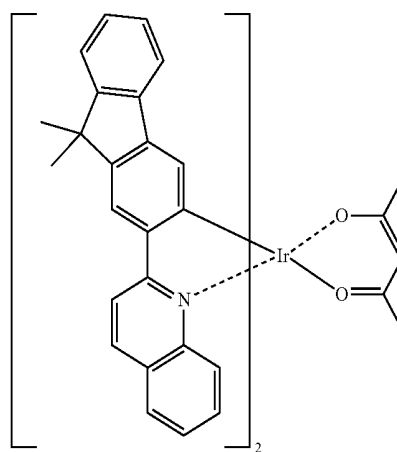
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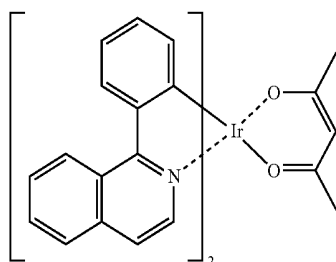
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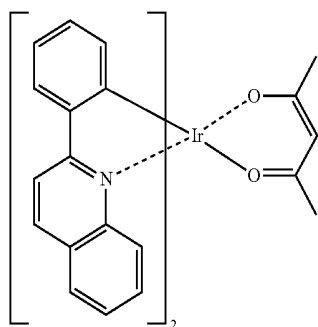
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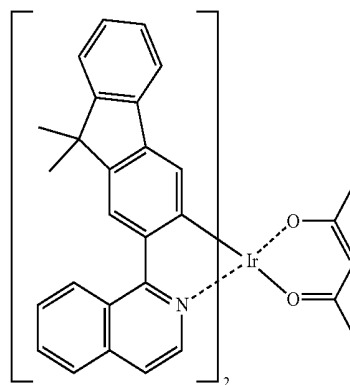
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PD16

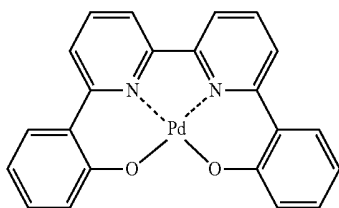


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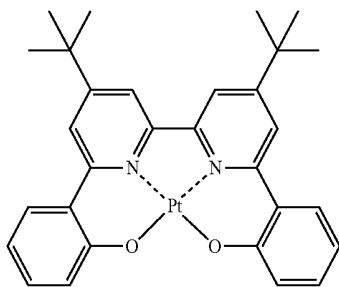


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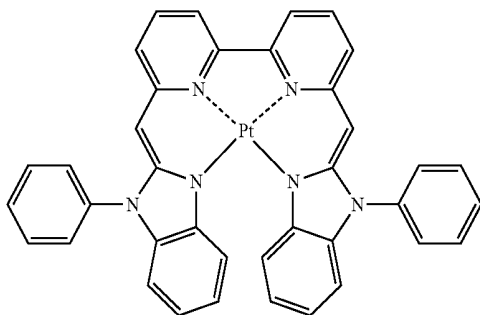
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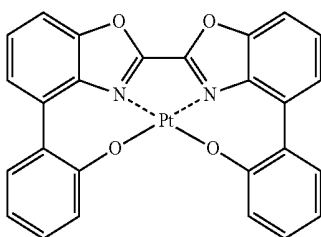
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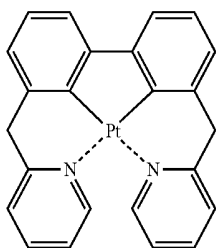
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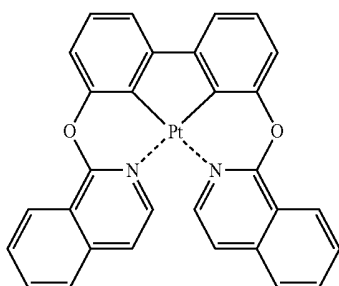
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PD25

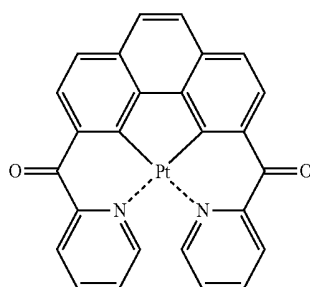


PD26

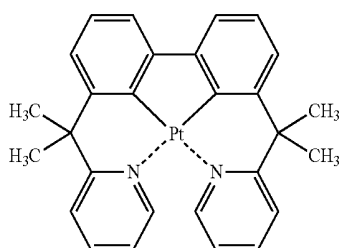


PD27

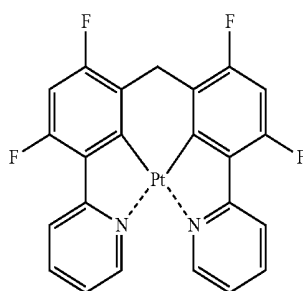
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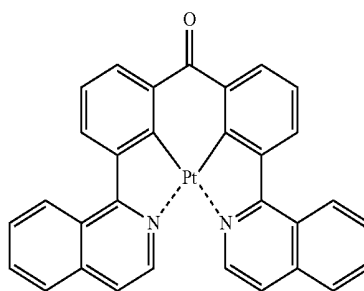
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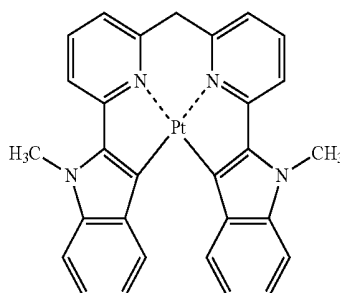
PD29



PD30

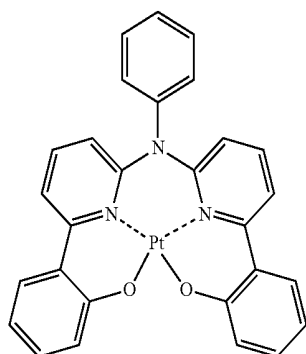
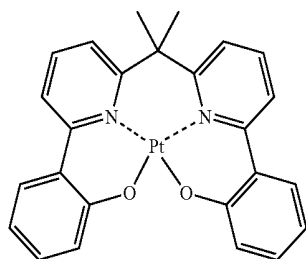
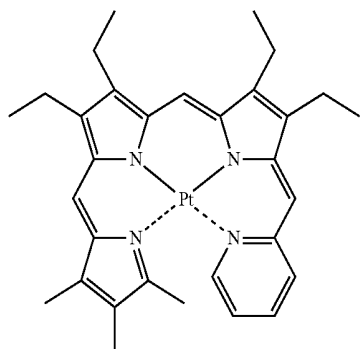
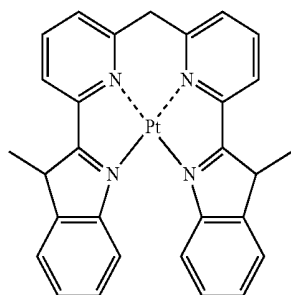
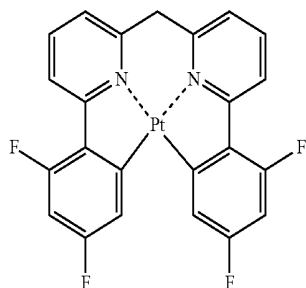


PD31



PD32

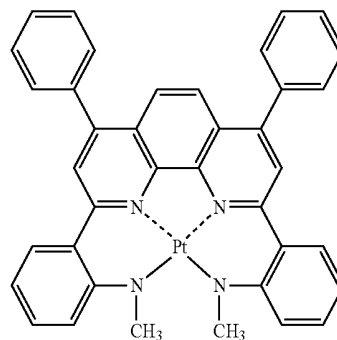
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PD33

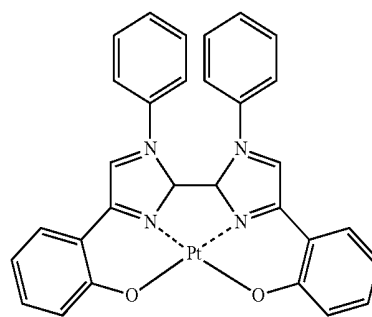
PD38



PD34

PD35

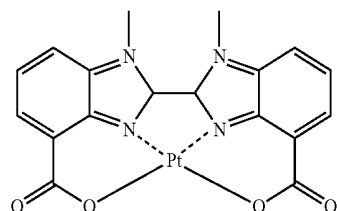
PD39



PD40

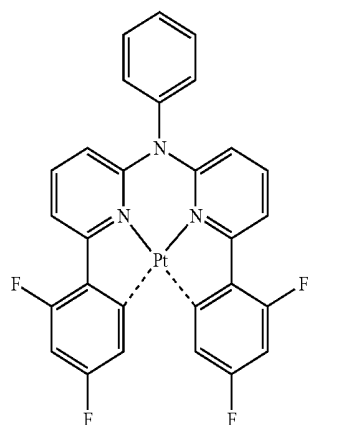
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PD41

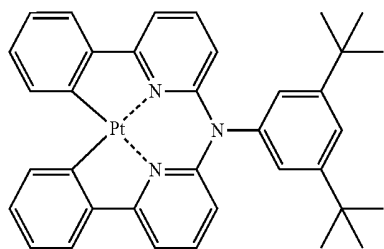
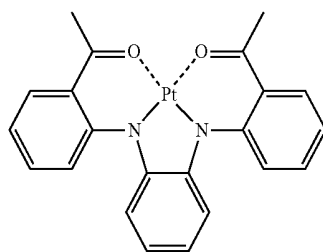
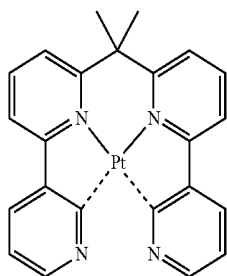
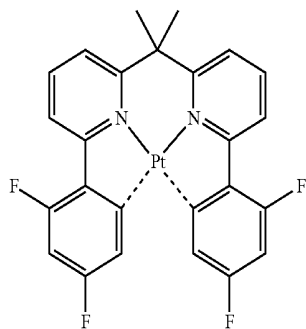
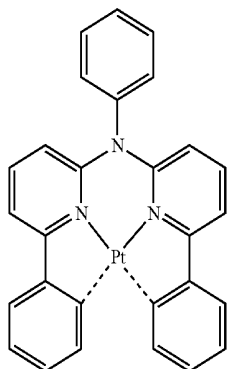


PD37

PD42

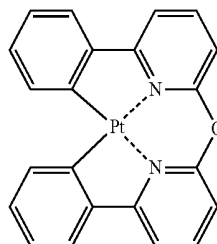


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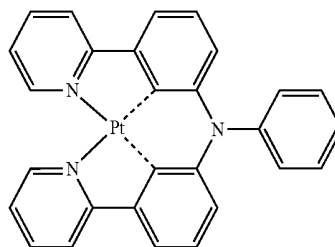
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PD43



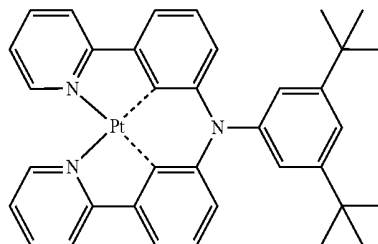
PD48

PD44



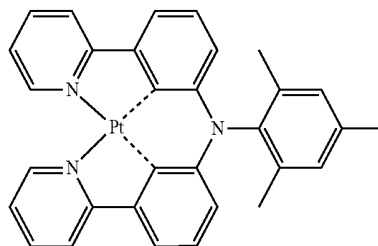
PD49

PD45



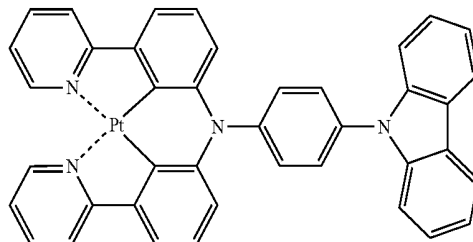
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PD46

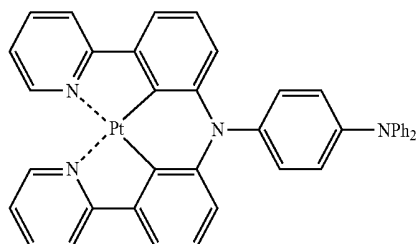


PD51

PD47

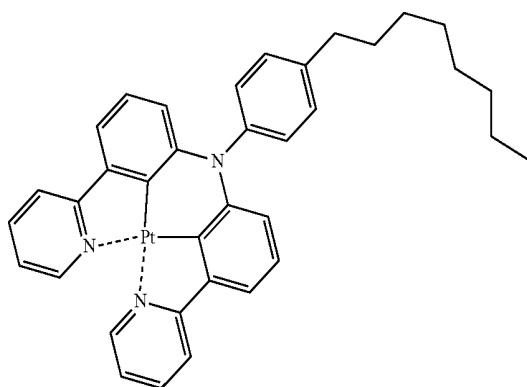
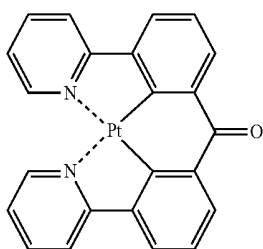
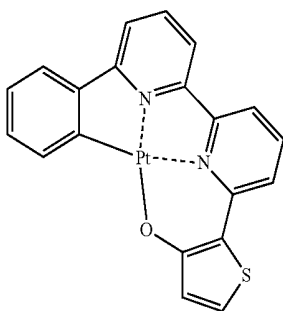
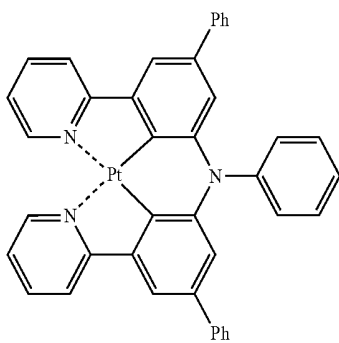
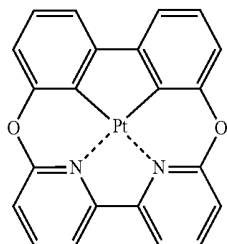


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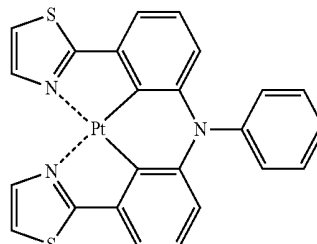


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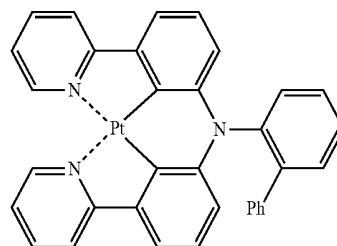
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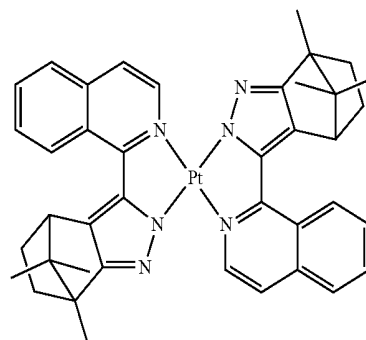
PD54



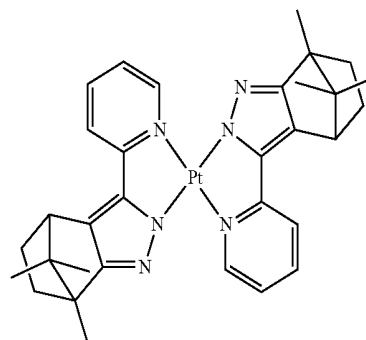
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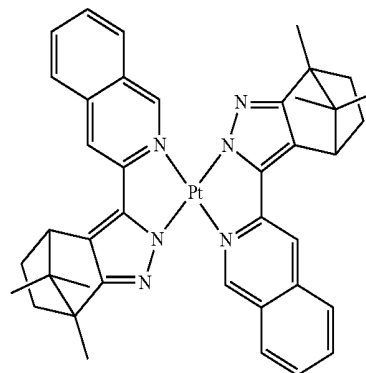
PD56



PD57



PD58



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PD59

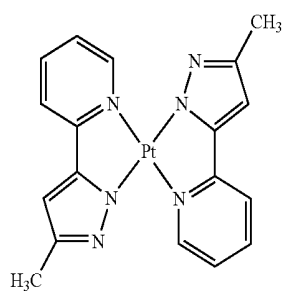
PD60

PD61

PD62

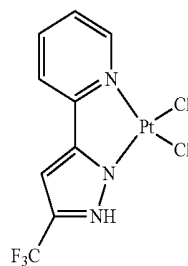
PD63

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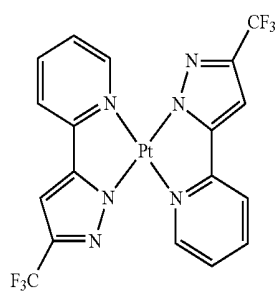


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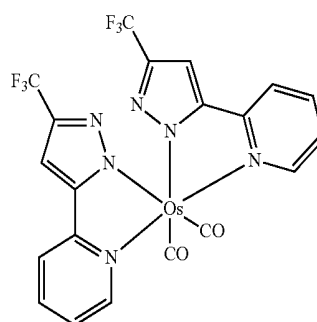
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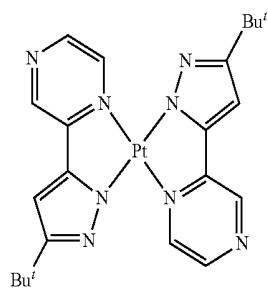
PD69



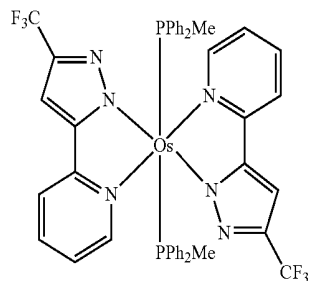
PD65



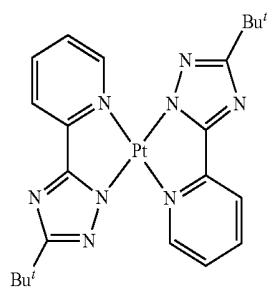
PD70



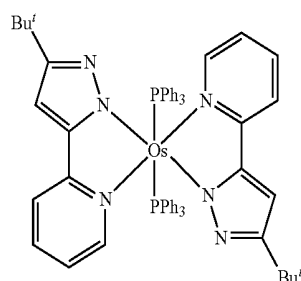
PD66



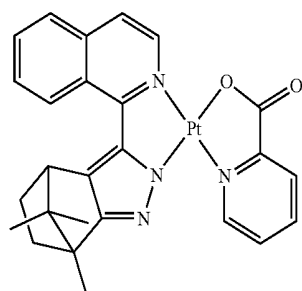
PD71



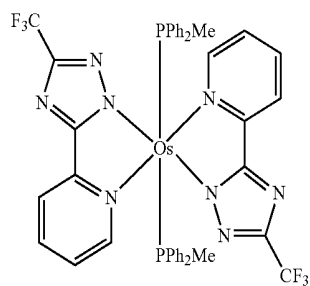
PD67



PD72

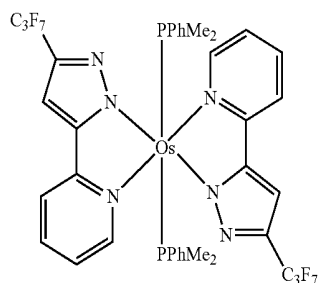


PD68

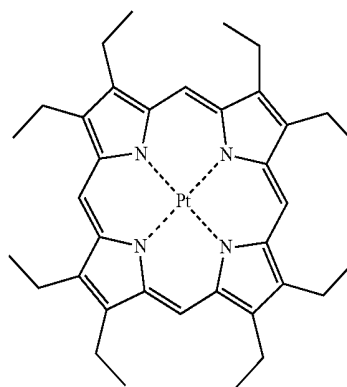


PD73

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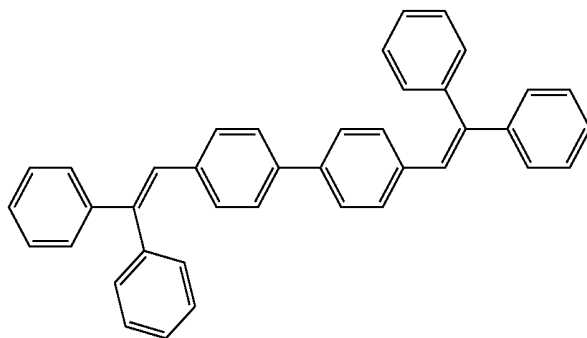
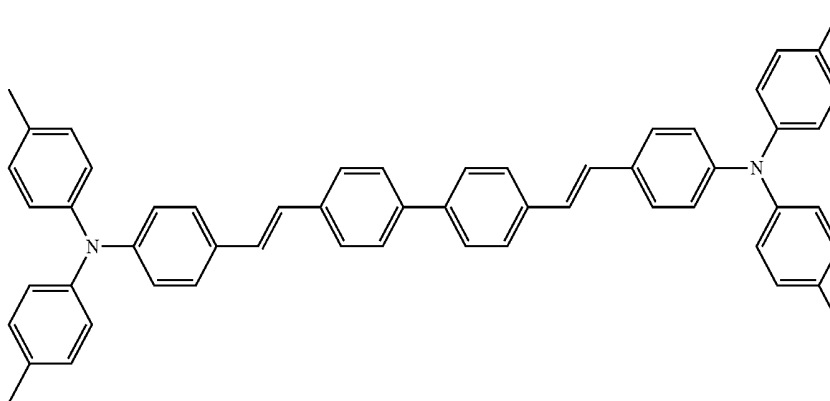
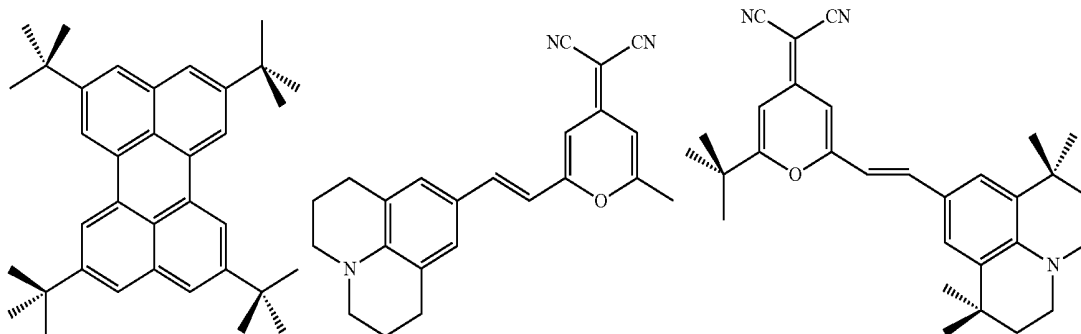
PD74



PtOEP

[0220] According to another embodiment, the phosphorescent dopant may include PtOEP.

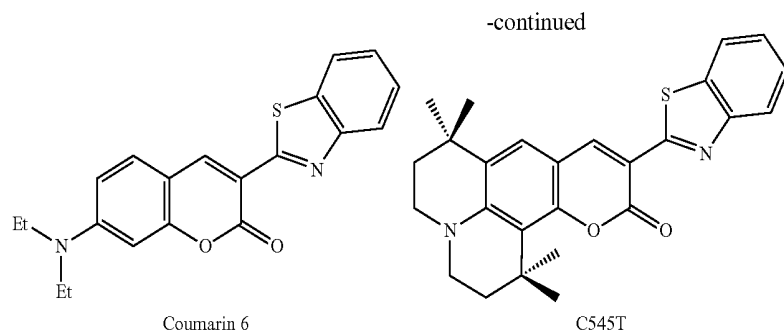
[0221] The fluorescent dopant may include at least one selected from DPAVB_i, BDAVB_i, TBPe, DCM, DCJT_B, Coumarin 6, and C545T.

DPVB_iDPAVB_i

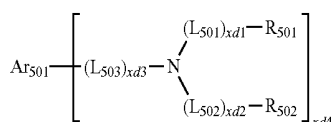
TBPe

DCM

DCJT_B



[0222] According to another embodiment, the fluorescent dopant may include a compound represented by Formula 501 below.



<Formula 501>

[0223] In Formula 501,

[0224] Ar_{501} may be selected from a naphthalene, a heptalene, a fluorene, a spiro-fluorene, a benzofluorene, a dibenzofluorene, a phenalene, a phenanthrene, an anthracene, a fluoranthene, a triphenylene, a pyrene, a chrysene, a naphthacene, a picene, a perylene, a pentaphene, and an indenoanthracene; and

[0225] a naphthalene group, a heptalene group, a fluorene group, a spiro-fluorene group, a benzofluorene group, a dibenzofluorene group, a phenalene group, a phenanthrene group, an anthracene group, a fluoranthene group, a triphenylene group, a pyrene group, a chrysene group, a naphthacene group, a picene group, a perylene group, a pentaphene group, and an indenoanthracene group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid group or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a $\text{C}_1\text{--C}_{60}$ alkyl group, a $\text{C}_2\text{--C}_{60}$ alkenyl group, a $\text{C}_2\text{--C}_{60}$ alkynyl group, a $\text{C}_1\text{--C}_{60}$ alkoxy, a $\text{C}_3\text{--C}_{10}$ cycloalkyl group, a $\text{C}_1\text{--C}_{10}$ heterocycloalkyl group, a $\text{C}_3\text{--C}_{10}$ cycloalkenyl group, a $\text{C}_1\text{--C}_{10}$ heterocycloalkenyl group, a $\text{C}_6\text{--C}_{60}$ aryl group, a $\text{C}_6\text{--C}_{60}$ aryloxy, a $\text{C}_6\text{--C}_{60}$ arylthio group, a $\text{C}_1\text{--C}_{60}$ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, and $-\text{Si}(\text{Q}_{501})(\text{Q}_{502})(\text{Q}_{503})$ (wherein Q_{501} to Q_{503} may be each independently selected from a hydrogen, a $\text{C}_1\text{--C}_{60}$ alkyl group, a $\text{C}_2\text{--C}_{60}$ alkenyl group, a $\text{C}_1\text{--C}_{60}$ aryl group, and a $\text{C}_1\text{--C}_{60}$ heteroaryl group);

[0226] L_{501} to L_{503} may be the same as explained in connection with L_{201} ;

[0227] R_{501} and R_{502} may be each independently selected from:

[0228] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a

pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a triazinyl group, a dibenzofuranyl group, and a dibenzothiophenyl group; and

[0229] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a triazinyl group, a dibenzofuranyl group, and a dibenzothiophenyl group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a $\text{C}_1\text{--C}_{20}$ alkyl group, a $\text{C}_1\text{--C}_{20}$ alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a triazinyl group, a dibenzofuranyl group, and a dibenzothiophenyl group;

[0230] x_{d1} to x_{d3} may be each independently selected from 0, 1, 2, and 3; and

[0231] x_{b4} may be selected from 1, 2, 3, and 4.

[0232] An amount of the dopant in the emission layer may be, e.g., in a range of about 0.01 part by weight to about 15 parts by weight, based on 100 parts by weight of the host, but is not limited thereto.

[0233] A thickness of the emission layer may be in a range of about 100 Å to about 1,000 Å, e.g., about 200 Å to about 600 Å. When the thickness of the emission layer is within this range, excellent light-emission characteristics may be obtained without a substantial increase in driving voltage.

[0234] Then, an electron transport region may be disposed on the emission layer.

[0235] The electron transport region may include at least one selected from a hole blocking layer, an electron transport layer (ETL), and an electron injection layer, but is not limited thereto.

[0236] For example, the electron transport region may have a structure of electron transport layer/electron injection layer or a structure of hole blocking layer/electron transport layer/electron injection layer, wherein layers of each structure are

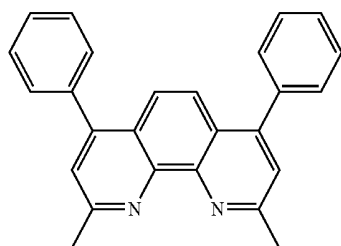
sequentially stacked from the emission layer in the stated order, but is not limited thereto.

[0237] According to an embodiment, the organic layer **150** of the organic light-emitting device includes an electron transport region disposed between the emission layer and the second electrode **190**, wherein the electron transport region includes the carbazole-based compound represented by Formula 1.

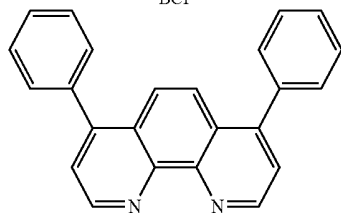
[0238] The electron transport region may include a hole blocking layer. The hole blocking layer may be formed, when the emission layer includes a phosphorescent dopant, to prevent diffusion of excitons or holes into an electron transport layer.

[0239] When the electron transport region includes a hole blocking layer, the hole blocking layer may be formed on the emission layer by using various methods, such as vacuum deposition, spin coating casting, a LB method, ink-jet printing, laser-printing, or laser-induced thermal imaging. When the hole blocking layer is formed by vacuum deposition or spin coating, deposition and coating conditions for the hole blocking layer may be determined by referring to the deposition and coating conditions for the hole injection layer.

[0240] The hole blocking layer may include, for example, at least one selected from BCP and Bphen, but is not limited thereto.



BCP



Bphen

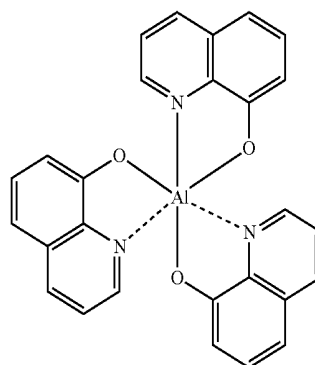
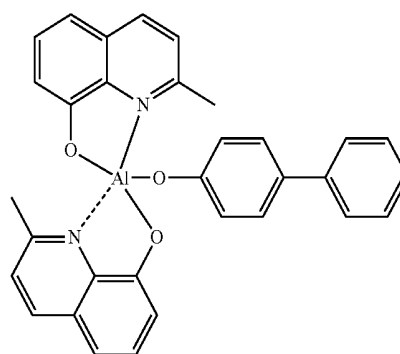
[0241] A thickness of the hole blocking layer may be in a range of about 20 Å to about 1000 Å, for example, about 30 Å to about 300 Å. When the thickness of the hole blocking layer is within this range, the hole blocking layer may have improved hole blocking ability without a substantial increase in driving voltage.

[0242] The electron transport region may include an electron transport layer. The electron transport layer may be formed on the emission layer or the hole blocking layer by using various methods, such as vacuum deposition, spin coating casting, a LB method, ink-jet printing, laser-printing, or laser-induced thermal imaging. When an electron transport layer is formed by vacuum deposition or spin coating, deposition and coating conditions for the electron transport layer may be determined by referring to the deposition and coating conditions for the hole injection layer.

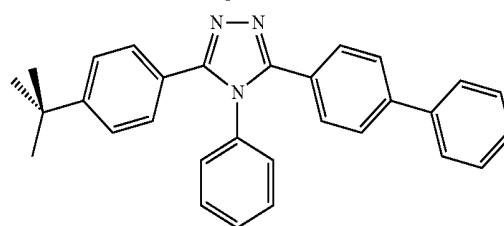
[0243] According to an embodiment, the organic layer **150** of the organic light-emitting device may include an electron

transport region disposed between the emission layer and the second electrode **190**. The electron transport region may include at least one selected from an electron transport layer and an electron injection layer.

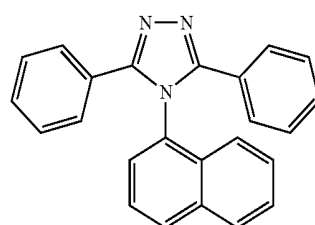
[0244] The electron transport layer may further include at least one selected from BCP, Bphen, Alq₃, Balq, TAZ, and NTAZ.

Alq₃

Balq

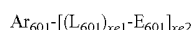


TAZ



NTAZ

[0245] According to another embodiment, the electron transport layer may include at least one compound selected from a compound represented by Formula 601 and a compound represented by Formula 602 illustrated below:



<Formula 601>

[0246] wherein in Formula 601,

[0247] Ar₆₀₁ may be selected from

[0248] a naphthalene, a heptalene, a fluorene, a spiro-fluorene, a benzofluorene, a dibenzofluorene, a phenalene, a phenanthrene, an anthracene, a fluoranthene, a triphenylene, a pyrene, a chrysene, naphthacene, a picene, a perylene, a pentaphene and an indenoanthracene;

[0249] a naphthalene, a heptalene, a fluorene, a spiro-fluorene, a benzofluorene, a dibenzofluorene, a phenalene, a phenanthrene, an anthracene, a fluoranthene, a triphenylene, a pyrene, a chrysene, a naphthacene, a picene, a perylene, a pentaphene, and an indenoanthracene, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid group or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, and —Si(Q₃₀₁)(Q₃₀₂)(Q₃₀₃) (wherein Q₃₀₁ to Q₃₀₃ may be each independently selected from a hydrogen, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₁-C₆₀ aryl group, and a C₁-C₆₀ heteroaryl group);

[0250] L₆₀₁ may be the same as explained in connection with L₂₀₁;

[0251] E₆₀₁ may be selected from

[0252] a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, a tetrazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a thiadiazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group; and

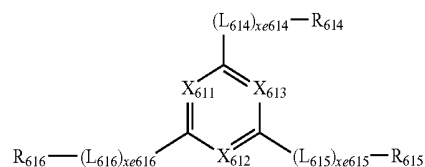
[0253] a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl

group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, a tetrazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a thiadiazolyl group, an imidazopyridinyl group, and an imidazopyrimidinyl group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, C₁-C₂₀ alkyl group, C₁-C₂₀ alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenyl group, a heptalenyl group, an indacenyl group, an acenaphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, a tetrazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a thiadiazolyl group, an imidazopyridinyl group, and an imidazopyrimidinyl group;

[0254] xe1 may be selected from 0, 1, 2, and 3; and

[0255] xe2 may be selected from 1, 2, 3, and 4.

<Formula 602>



[0256] wherein in Formula 602,

[0257] X_{611} may be N or C-(L_{611}) $_{xe611}$ - R_{611} , X_{612} may be N or C-(L_{612}) $_{xe612}$ - R_{612} , X_{613} may be N or C-(L_{613}) $_{xe613}$ - R_{613} and, at least one selected from X_{611} to X_{613} may be N;

[0258] L_{611} to L_{616} may be the same as explained in connection with L_{201} ;

[0259] R_{611} to R_{616} are each independently selected from

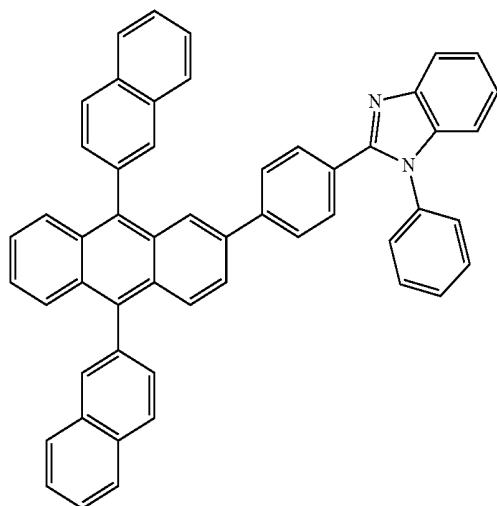
[0260] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group and a triazinyl group; and

[0261] a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, and a triazinyl group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, and a triazinyl group; and

[0262] $xe611$ to $xe616$ may be each independently selected from 0, 1, 2, and 3.

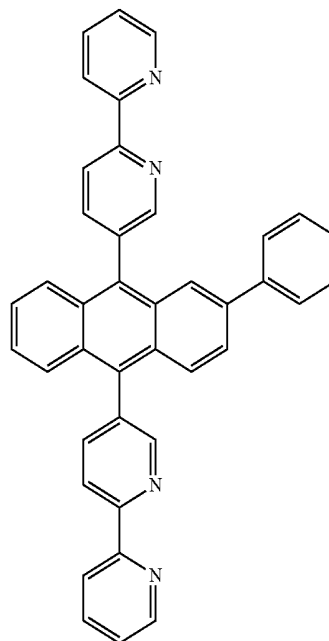
[0263] The compound represented by Formula 601 and the compound represented by Formula 602 may each be selected from Compounds ET1 to ET15 illustrated below.

ET1

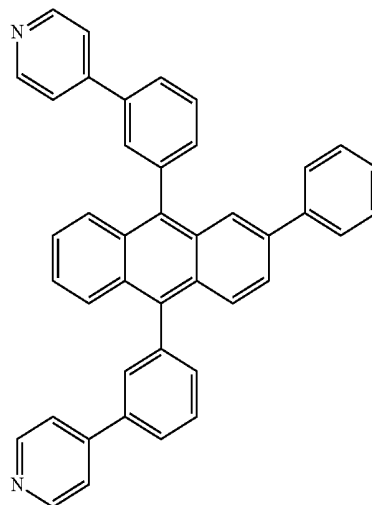


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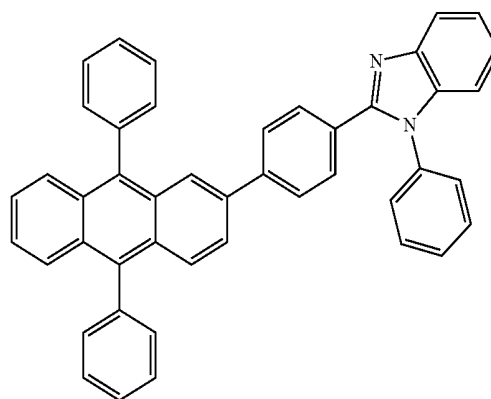
ET2



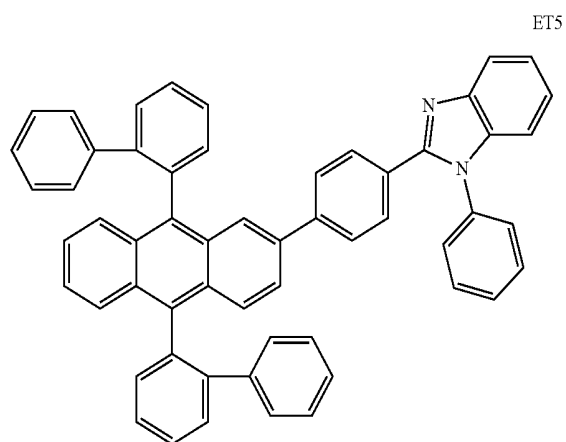
ET3



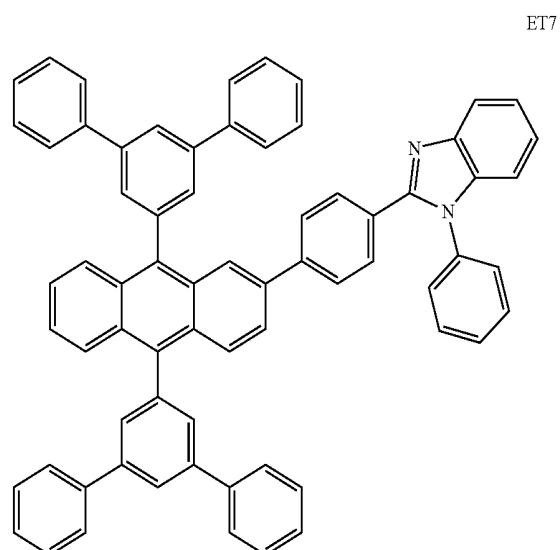
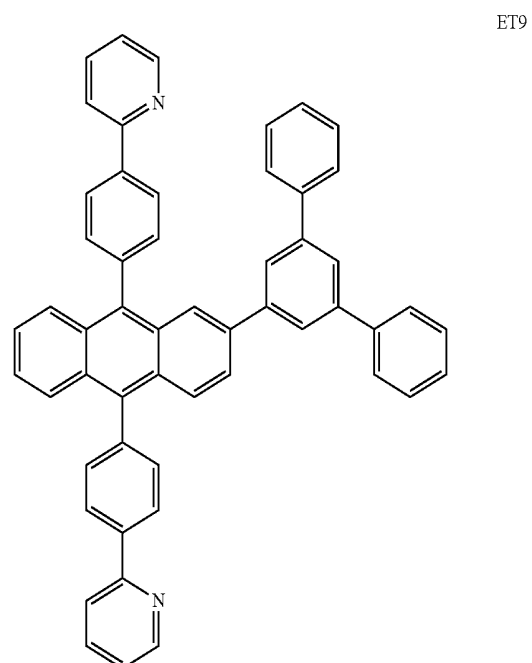
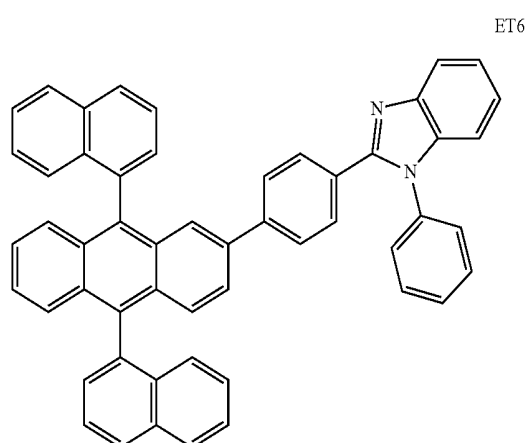
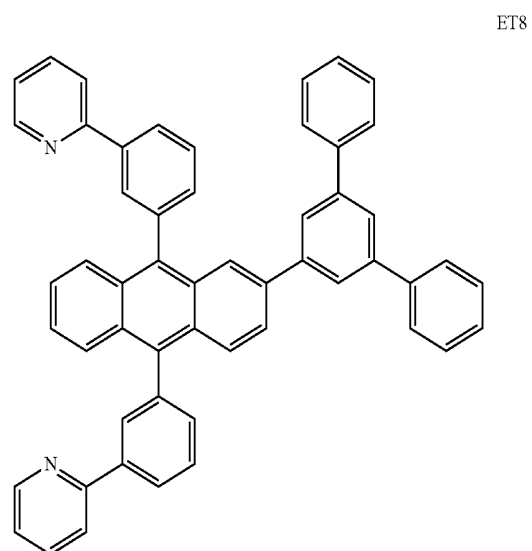
ET4



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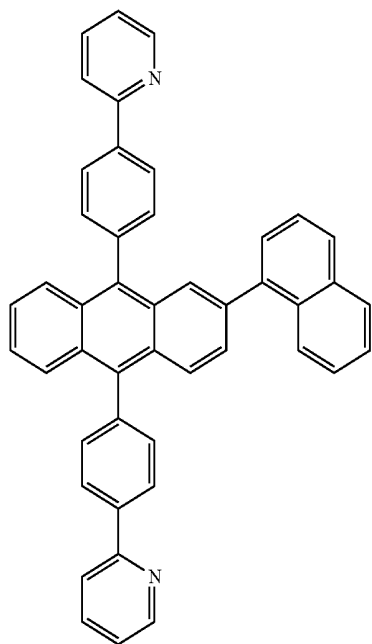


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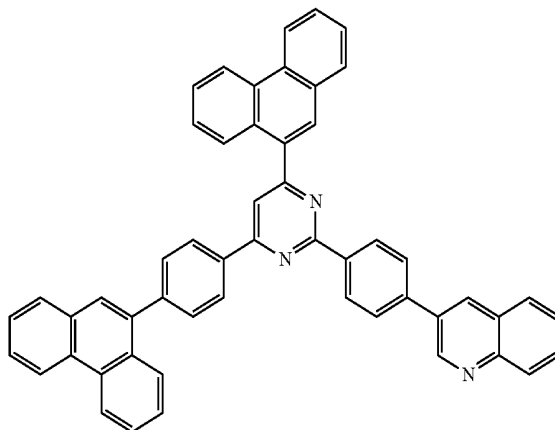
-continued

ET10



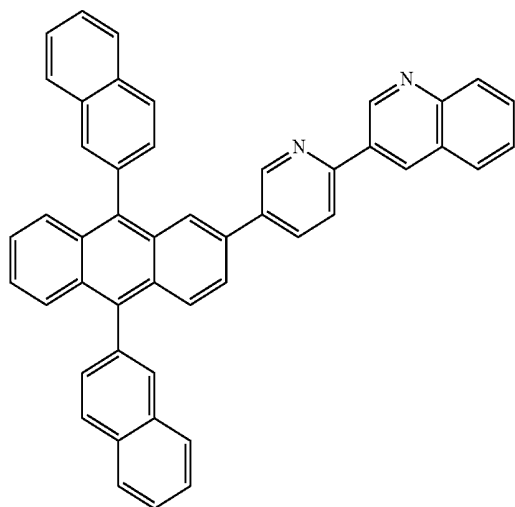
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ET13

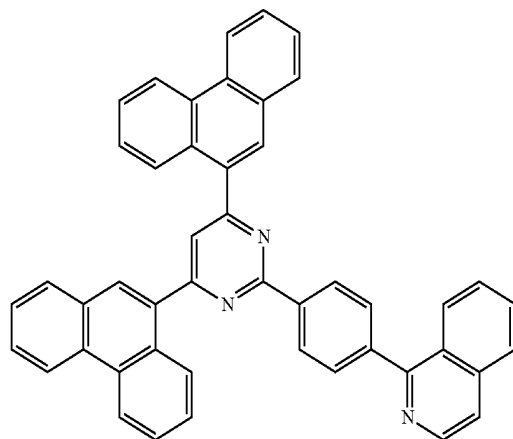


ET14

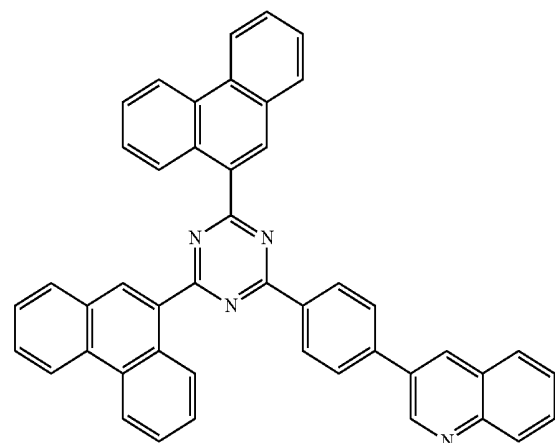
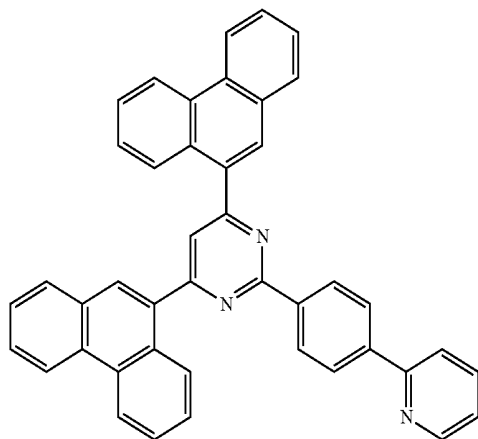
ET11



ET15



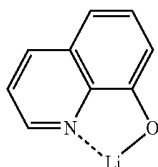
ET12



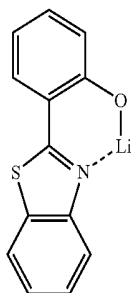
[0264] A thickness of the electron transport layer may be in a range of about 100 Å to about 1,000 Å, e.g., about 150 Å to about 500 Å. When the thickness of the electron transport layer is within the range, the electron transport layer may have satisfactory electron transport characteristics without a substantial increase in driving voltage.

[0265] Also, the electron transport layer may further include, in addition to the materials described above, a metal-containing material.

[0266] The metal-containing material may include a Li complex. The Li complex may include, for example, Compound ET-D1 (lithium quinolate, LiQ) or ET-D2.



ET-D1



ET-D2

[0267] The electron transport region may include an electron injection layer that allows electrons to be easily provided from the second electrode **190**.

[0268] The electron injection layer may be formed on the electron transport layer by using various methods, such as vacuum deposition, spin coating casting, a LB method, ink-jet printing, laser-printing, or laser-induced thermal imaging. When an electron injection layer is formed by vacuum deposition or spin coating, deposition and coating conditions for the electron injection layer may be determined by referring to the deposition and coating conditions for the hole injection layer.

[0269] The electron injection layer may include at least one selected from, LiF, NaCl, CsF, Li₂O, BaO, and LiQ.

[0270] A thickness of the electron injection layer may be in a range of about 1 Å to about 100 Å, e.g., about 3 Å to about 90 Å. When the thickness of the electron injection layer is within the range, the electron injection layer may have satisfactory electron injection characteristics without a substantial increase in driving voltage.

[0271] The second electrode **190** may be disposed on the organic layer **150** having such a structure. The second electrode **190** may be a cathode that is an electron injection electrode, and in this regard, a material for forming the second electrode **190** may be a material having a low work function, and such a material may be metal, alloy, an electrically conductive compound, or a mixture thereof. Detailed examples of the second electrode **190** are lithium (Li), magnesium (Mg), aluminum (Al), aluminum-lithium (Al—Li), calcium (Ca), magnesium-indium (Mg—In), or magnesium-silver (Mg—Ag). According to another embodiment, the material for forming the second electrode **190** may be ITO or IZO. The second electrode **190** may be a reflective electrode, a semi-transmissive electrode, or a transmissive electrode.

[0272] Hereinbefore, the organic light-emitting device has been described with reference to FIG. 1, but is not limited thereto.

[0273] A C₁-C₆₀ alkyl group used herein refers to a linear or branched aliphatic hydrocarbon monovalent group having 1 to 60 carbon atoms, and detailed examples thereof are a

methyl group, an ethyl group, a propyl group, an isobutyl group, a sec-butyl group, a tert-butyl group, a pentyl group, an iso-amyl group, and a hexyl group. A C₁-C₆₀ alkylene group used herein refers to a divalent group having the same structure as the C₁-C₆₀ alkyl group.

[0274] A C₁-C₆₀ alkoxy group used herein refers to a monovalent group represented by —OA₁₀₁ (wherein A₁₀₁ is the C₁-C₆₀ alkyl group), and detailed examples thereof are a methoxy group, an ethoxy group, and an isopropoxy group.

[0275] A C₂-C₆₀ alkenyl group used herein refers to a hydrocarbon group formed by substituting at least one carbon double bond in the middle or terminal of the C₂-C₆₀ alkyl group, and detailed examples thereof are an ethenyl group, a propenyl group, and a butenyl group. A C₂-C₆₀ alkylene group used herein refers to a divalent group having the same structure as the C₂-C₆₀ alkyl group.

[0276] A C₂-C₆₀ alkynyl group used herein refers to a hydrocarbon group formed by substituting at least one carbon triple bond in the middle or terminal of the C₂-C₆₀ alkyl group, and detailed examples thereof are an ethynyl group, and a propynyl group. A C₂-C₆₀ alkylene group used herein refers to a divalent group having the same structure as the C₂-C₆₀ alkyl group.

[0277] A C₃-C₁₀ cycloalkyl group used herein refers to a monovalent hydrocarbon monocyclic group having 3 to 10 carbon atoms, and detailed examples thereof are a cyclopropyl group, a cyclobutyl group, a cyclopentyl group, a cyclohexyl group, and a cycloheptyl group. A C₃-C₁₀ cycloalkylene group used herein refers to a divalent group having the same structure as the C₃-C₁₀ cycloalkyl group.

[0278] A C₁-C₁₀ heterocycloalkyl group used herein refers to a monovalent monocyclic group having at least one hetero atom selected from N, O, P, and S as a ring-forming atom and 1 to 10 carbon atoms, and detailed examples thereof are a tetrahydrofuranyl group, and a tetrahydrothiophenyl group. A C₁-C₁₀ heterocycloalkylene group used herein refers to a divalent group having the same structure as the C₁-C₁₀ heterocycloalkyl group.

[0279] A C₃-C₁₀ cycloalkenyl group used herein refers to a monovalent monocyclic group that has 3 to 10 carbon atoms and at least one double bond in the ring thereof and does not have aromaticity, and detailed examples thereof are a cyclopentenyl group, a cyclohexenyl group, and a cycloheptenyl group. A C₃-C₁₀ cycloalkenylene group used herein refers to a divalent group having the same structure as the C₃-C₁₀ cycloalkenyl group.

[0280] A C₁-C₁₀ heterocycloalkenyl group used herein refers to a monovalent monocyclic group that has at least one hetero atom selected from N, O, P, and S as a ring-forming atom, 1 to 10 carbon atoms, and at least one double bond in its ring. Detailed examples of the C₁-C₁₀ heterocycloalkenyl group are a 2,3-dihydrofuranyl group and a 2,3-dihydrothiophenyl group. A C₁-C₁₀ heterocycloalkenylene group used herein refers to a divalent group having the same structure as the C₁-C₁₀ heterocycloalkenyl group.

[0281] A C₆-C₆₀ aryl group used herein refers to a monovalent group having a carbocyclic aromatic system having 6 to 60 carbon atoms, and a C₆-C₆₀ arylene group used herein refers to a divalent group having a carbocyclic aromatic system having 6 to 60 carbon atoms. Detailed examples of the C₆-C₆₀ aryl group are a phenyl group, a naphthyl group, an anthracenyl group, a phenanthrenyl group, a pyrenyl group, and a chrysenyl group. When the C₆-C₆₀ aryl group and the

C₆-C₆₀ arylene group each include two or more rings, the rings may be fused to each other.

[0282] A C₁-C₆₀ heteroaryl group used herein refers to a monovalent group having a carbocyclic aromatic system that has at least one hetero atom selected from N, O, P, and S as a ring-forming atom, and 1 to 60 carbon atoms. A C₁-C₆₀ heteroarylene group used herein refers to a divalent group having a carbocyclic aromatic system that has at least one hetero atom selected from N, O, P, and S as a ring-forming atom, and 1 to 60 carbon atoms. Examples of the C₁-C₆₀ heteroaryl group are a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group. When the C₁-C₆₀ heteroaryl group and the C₁-C₆₀ heteroarylene group each include two or more rings, the rings may be fused to each other.

[0283] A C₆-C₆₀ aryloxy group used herein indicates —OA₁₀₂ (wherein A₁₀₂ is the C₆-C₆₀ aryl group), and a C₆-C₆₀ arylthio group used herein indicates —SA₁₀₃ (wherein A₁₀₃ is the C₆-C₆₀ aryl group).

[0284] A monovalent non-aromatic condensed polycyclic group used herein refers to a monovalent group (for example, having 8 to 60 carbon atoms) that has two or more rings condensed to each other, only carbon atoms as a ring forming atom, and non-aromaticity in the entire molecular structure. A detailed example of the monovalent non-aromatic condensed polycyclic group is a fluorenyl group. A divalent non-aromatic condensed polycyclic group used herein refers to a divalent group having the same structure as the monovalent non-aromatic condensed polycyclic group.

[0285] A monovalent non-aromatic condensed heteropolycyclic group used herein refers to a monovalent group (for example, having 1 to 60 carbon atoms) that has two or more rings condensed to each other, has a heteroatom selected from N, O, P, and S, other than carbon atoms, as a ring forming atom, and has non-aromaticity in the entire molecular structure. An example of the monovalent non-aromatic condensed heteropolycyclic group is a carbazolyl group. A divalent non-aromatic condensed heteropolycyclic group used herein refers to a divalent group having the same structure as the monovalent non-aromatic condensed heteropolycyclic group.

[0286] At least one of substituents of the substituted carbazolyl group, substituted C₃-C₁₀ cycloalkylene group, substituted C₁-C₁₀ heterocycloalkylene group, substituted C₃-C₁₀ cycloalkenylene group, substituted C₁-C₁₀ heterocycloalkenylene group, substituted C₆-C₆₀ arylene group, substituted C₁-C₆₀ heteroarylene group, substituted divalent non-aromatic condensed polycyclic group, substituted divalent non-aromatic condensed heteropolycyclic group, substituted C₁-C₆₀ alkyl group, substituted C₂-C₆₀ alkenyl group, substituted C₂-C₆₀ alkynyl group, substituted C₁-C₆₀ alkoxy group, substituted C₃-C₁₀ cycloalkyl group, substituted C₁-C₁₀ heterocycloalkyl group, substituted C₃-C₁₀ cycloalkenyl group, substituted C₁-C₁₀ heterocycloalkenyl group, substituted C₆-C₆₀ aryl group, substituted C₆-C₆₀ aryloxy group, substituted C₆-C₆₀ arylthio group, substituted C₁-C₆₀ heteroaryl group, substituted monovalent non-aromatic condensed polycyclic group, and substituted monovalent non-aromatic condensed heteropolycyclic group may be selected from

[0287] a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a

phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group;

[0288] a C₁-C₆₀ alkyl group, a C₁-C₆₀ alkenyl group, a C₁-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid group or a salt thereof, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, —Si(Q₁₁)(Q₁₂)(Q₁₃), —B(Q₁₄)(Q₁₅), and —N(Q₁₆)(Q₁₇);

[0289] a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group;

[0290] a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic heterocondensed polycyclic group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid group or a salt thereof, a C₁-C₆₀ alkyl group, a C₁-C₆₀ alkenyl group, a C₁-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, —Si(Q₂₁)(Q₂₂), —B(Q₂₃)(Q₂₄)(Q₂₅), and —N(Q₂₆)(Q₂₇); and

[0291] —Si(Q₃₁)(Q₃₂)(Q₃₃), —B(Q₃₄)(Q₃₅) and —N(Q₃₆)(Q₃₇);

[0292] wherein Q₁₁ to Q₁₇, Q₂₁ to Q₂₇, and Q₃₁ to Q₃₇ may be each independently selected from a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group.

[0293] For example, at least one substituent selected from substituents of the substituted carbazolyl group, the substituted C₃-C₁₀ cycloalkylene group, substituted C₁-C₁₀ hetero-

cycloalkylene group, substituted C_3 - C_{10} cycloalkenylene group, substituted C_1 - C_{10} heterocycloalkenylene group, substituted C_6 - C_{60} arylene group, substituted C_1 - C_{60} heteroarylene group, substituted divalent non-aromatic condensed polycyclic group, substituted divalent non-aromatic condensed heteropolycyclic group, substituted C_1 - C_{60} alkyl group, substituted C_2 - C_{60} alkenyl group, substituted C_2 - C_{60} alkynyl group, substituted C_1 - C_{60} alkoxy group, substituted C_3 - C_{10} cycloalkyl group, substituted C_1 - C_{10} heterocycloalkyl group, substituted C_3 - C_{10} cycloalkenyl group, substituted C_1 - C_{10} heterocycloalkenyl group, substituted C_6 - C_{60} aryl group, substituted C_6 - C_{60} aryloxy group, substituted C_6 - C_{60} arylthio group, substituted C_1 - C_{60} heteroaryl group, substituted monovalent non-aromatic condensed polycyclic group, and substituted monovalent non-aromatic condensed heteropolycyclic group may be selected from

[0294] a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group;

[0295] a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenyl group, a heptalenyl group, an indacenyl group, an acenaphthyl group, a fluorenyl group, a spirofluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isooxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzooxazolyl group, an isobenzooxazolyl group, a triazolyl group, a tetrazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a thiadiazolyl group, —Si(Q₁)(Q₁₂)(Q₁₃), —B(Q₁₄)(Q₁₅), and —N(Q₁₆)(Q₁₇);

[0296] a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenyl group, a heptalenyl group, an indace-

nyl group, an acenaphthyl group, a fluorenyl group, a spirofluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isooxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzooxazolyl group, an isobenzooxazolyl group, a triazolyl group, a tetrazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group;

[0297] a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenyl group, a heptalenyl group, an indacenyl group, an acenaphthyl group, a fluorenyl group, a spirofluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isooxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzooxazolyl group, an isobenzooxazolyl group, a triazolyl group, a tetrazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, a C_1 - C_{60} alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenyl group, a heptalenyl group, an indace-

indacenyl group, an acenaphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinoliny group, an isoquinoliny group, a benzoquinoliny group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnoliny group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenaziny group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, a tetrazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a thiadiazolyl group, $-\text{Si}(\text{Q}_{21})(\text{Q}_{22})(\text{Q}_{23})$, $-\text{B}(\text{Q}_{24})(\text{Q}_{25})$, and $-\text{N}(\text{Q}_{26})(\text{Q}_{27})$; and

[0298] $-\text{Si}(\text{Q}_{31})(\text{Q}_{32})(\text{Q}_{33})$, $-\text{B}(\text{Q}_{34})(\text{Q}_{35})$ and $-\text{N}(\text{Q}_{36})(\text{Q}_{37})$;

[0299] wherein Q_{11} to Q_{17} , Q_{21} to Q_{27} , and Q_{31} to Q_{37} may be each independently selected from a hydrogen, a deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a $\text{C}_1\text{-C}_{60}$ alkyl group, a $\text{C}_2\text{-C}_{60}$ alkenyl group, a $\text{C}_2\text{-C}_{60}$ alkynyl group, a $\text{C}_1\text{-C}_{60}$ alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenyl group, a heptalenyl group, an indacenyl group, an acenaphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinoliny group, an isoquinoliny group, a benzoquinoliny group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnoliny group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenaziny group, a benzoimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, a tetrazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group.

[0300] The term “Ph” used herein refers to a phenyl group, the term “Me” used herein refers to a methyl group, the term “Et” used herein refers to an ethyl group, and the term “ter-Bu” or “Bu^t” used herein refers to a tert-butyl group.

[0301] Hereinafter, an organic light-emitting device according to an embodiment will be described in detail with reference to Synthesis Examples and Examples. The wording “B was used instead of A” used in describing Synthesis Examples means that a molar equivalent of A was identical to a molar equivalent of B.

[0302] The following Examples and Comparative Examples are provided in order to highlight characteristics of one or more embodiments, but it will be understood that the Examples and Comparative Examples are not to be construed as limiting the scope of the embodiments, nor are the Comparative Examples to be construed as being outside the scope of the embodiments. Further, it will be understood that the embodiments are not limited to the particular details described in the Examples and Comparative Examples.

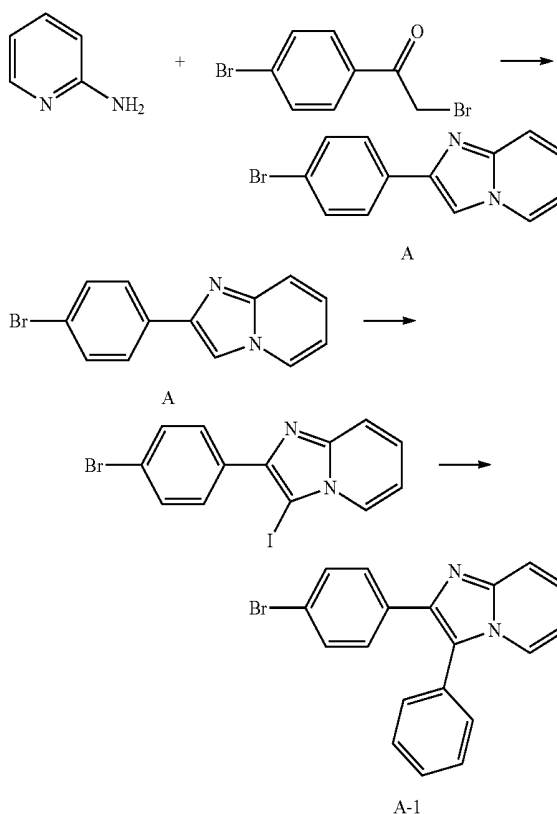
EXAMPLES

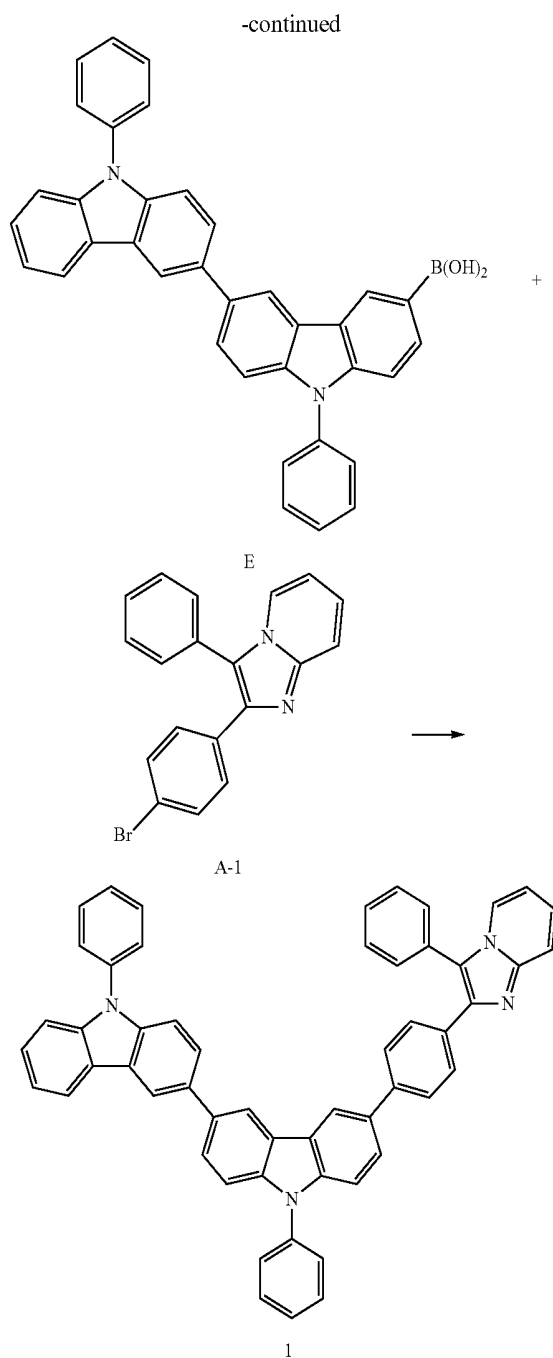
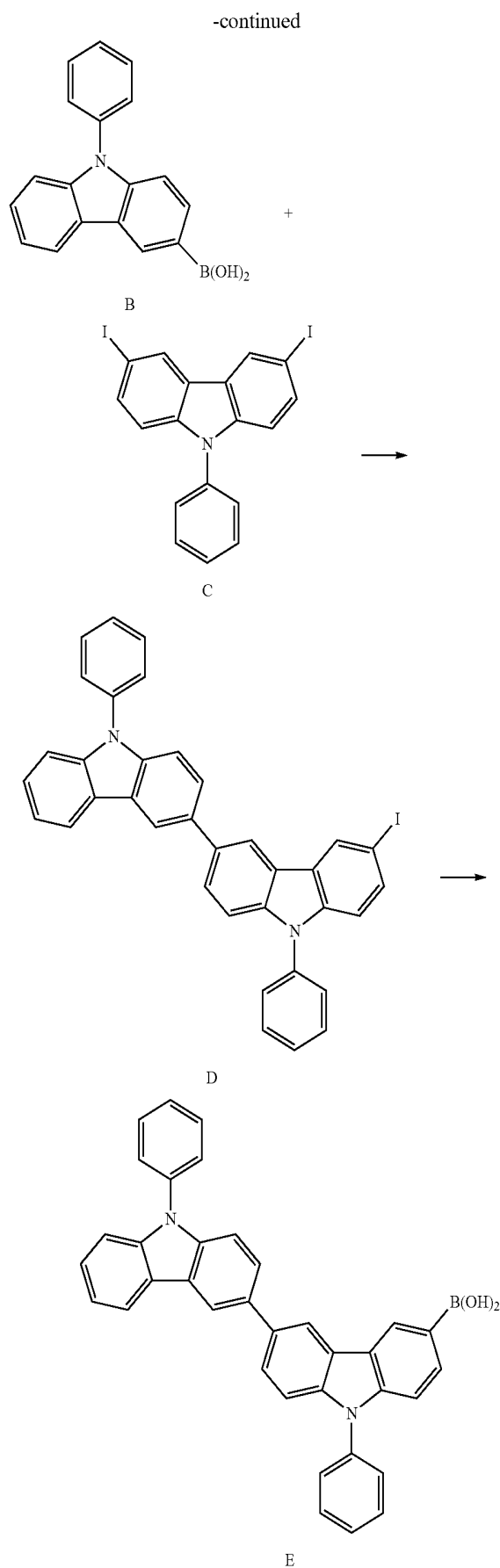
Synthesis Example

Synthesis Example 1

Synthesis of Compound 1

[0303]





[0304] Synthesis of Intermediate A

[0305] 3.39 g of 2-amino pyridine and 10 g of 2,4-dibromoacetophenone were dissolved in 150 ml of ethanol and the result was refluxed for 12 hours. Then, the result was cooled to ambient temperature to produce a white solid, and the white solid was washed with a aqueous NaHCO_3 saturated solution. The result was recrystallized using a dichloromethane/hexane solution to obtain 8.02 g (yield of 82%) of Intermediate A.

[0306] Synthesis of Intermediate A-1

[0307] 5 g (18.3 mmol) of Intermediate A and 4.12 g (18.3 mmol) of N-iodosuccinimide (NIS) were dissolved in an acetonitrile solvent and the result was stirred at ambient temperature for an hour, and then 100 ml of chloroform was added thereto. The result was washed with 10% aqueous

sodium hydroxide solution, and then washed with aqueous sodium thiosulfate saturated solution and water. After that, the result was dried with anhydrous magnesium sulfate and filtered, and then the solvent was removed. The obtained solid was washed with methanol and filtered to obtain 5.8 g (yield of 79%) of an iodine compound. The obtained iodine compound and 1.76 g (14.5 mmol) of a phenylboronic acid were added to a mixed solvent of aqueous potassium carbonate solution and tetrahydrofuran (THF). 335 mg of $\text{Pd}(\text{PPh}_3)_4$ was added to the result while stirring the result, and then the result was refluxed for 24 hours. The reaction solution was extracted with dichloromethane, the residual water was removed from the organic layer by using anhydrous magnesium sulfate, and then the result was dried under reduced pressure. After that, the result was separated by column chromatography (ethyl acetate:n-hexane) to obtain 2.93 g (yield of 58%) of Intermediate A-1.

[0308] Synthesis of Intermediate E

[0309] 8.6 g (30 mmol) of Compound B and 22.3 g (45 mmol) of Compound C were dissolved in a mixed solvent of 20.7 g (150 mmol) of aqueous potassium carbonate solution and THF. 1.7 g (1.5 mmol) of $\text{Pd}(\text{PPh}_3)_4$ was added to the result while stirring the result, and then the result was refluxed for 8 hours. The result was cooled to ambient temperature to obtain a solid, and then the obtained solid was washed with water, ethanol, and THF and filtered, thereby obtaining Compound D. 19.9 g (32.6 mmol) of Compound D obtained therein was dissolved in THF, and then 17.4 ml (39.1 mmol) of butyllithium was added dropwise thereto at a temperature of -78°C ., and after an hour, 3.8 ml (50.7 mmol) of trimethyl borate was added thereto. The result was heated to ambient temperature, and after an hour, 2N aqueous hydrochloric acid solution was added thereto and the result was stirred for 3 hours. The solid obtained was washed with toluene and filtered, thereby obtaining 13.4 g (yield of 78%) of Intermediate E.

[0310] Synthesis of Compound 1

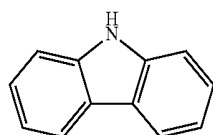
[0311] 10.5 g (30.0 mmol) of Intermediate A-1 and 15.9 g (30.0 mmol) of Intermediate E were dissolved in a mixed solvent of 20 g (150.2 mmol) of aqueous K_2CO_3 solution and THF. 1.7 g (1.51 mmol) of $\text{Pd}(\text{PPh}_3)_4$ was added to the result while stirring the result, and then the result was refluxed for 10 hours. The result was cooled to ambient temperature, and the solid obtained was washed with water, ethanol and THF and filtered, thereby obtaining 18.5 g (yield of 82%) of Compound 1.

[0312] ^1H NMR (300 MHz, CDCl_3) 8.49 (2H), 8.32 (3H), 8.15 (1H), 8.01-7.95 (3H), 7.86-7.81 (7H), 7.79 (2H), 7.64-7.56 (6H), 7.53-7.42 (8H), 7.37 (1H), 7.26-7.19 (2H), 6.98 (1H)

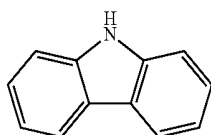
Synthesis Example 2

Synthesis of Compound 10

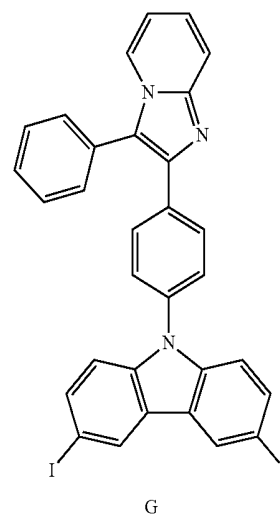
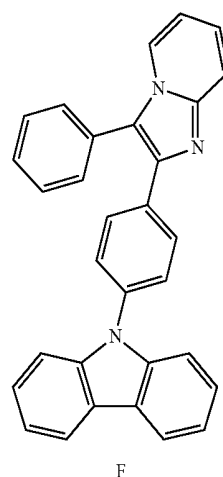
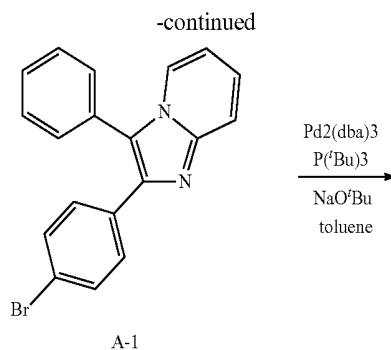
[0313]

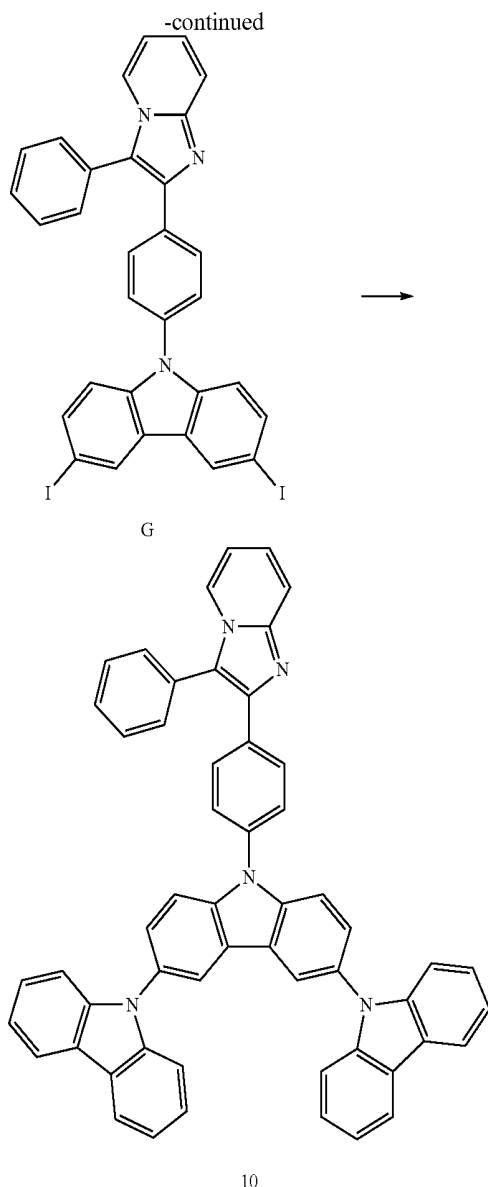


+



+



**[0314]** Synthesis of Intermediate G

[0315] 9.6 g (48 mmol) of carbazole, 14.0 g (40 mmol) of Intermediate A-1, 0.72 g (0.8 mmol) of Tris(dibenzylideneacetone)dipalladium(0):Pd₂(dba)₃, 160 mg (1.6 mmol) of P(t-Bu)₃ and 5.76 g (60.0 mmol) of t-BuONa were dissolved in toluene and the result was stirred at a temperature of 85° C. for 3 hours. Then, the result was cooled to ambient temperature and an extraction process was performed three times by using distilled water and ethyl acetate. After that, the extracted organic layers were dried by using magnesium sulfate, and the residue obtained after evaporating the solvent was separation-purified by silicagel column chromatography, thereby obtaining Intermediate F. 15.3 g (35.2 mmol) of Intermediate F obtained therein was completely dissolved in dichloromethane, and then 5.1 g (20.1 mmol) of iodine and 3.4 g (15.9 mmol) of KIO₃ were respectively divided into 5 portions, which then were sequentially added to the result. The result was stirred for 8 hours and then washed with methanol to obtain 12.8 g (yield of 53%) Intermediate G. The

obtained compound was confirmed by LC-MS (M+ 687.8) and sublimation-purified to have an HPLC purity of 99.1%.

[0316] Synthesis of Compound 10

[0317] Compound 10 (yield of 77%) was obtained by using the same coupling method as used in synthesizing Intermediate F with carbazole and Intermediate G.

[0318] ¹H NMR (300 MHz, CDCl₃) 8.53 (2H), 8.42 (1H), 8.15-8.09 (3H), 7.95-7.90 (6H), 7.86-7.79 (4H), 7.69 (2H), 7.63-7.48 (8H), 7.39-7.34 (3H), 7.26-7.18 (5H), 7.01 (1H)

Synthesis Example 3

Synthesis of Compound 19

[0319] Compound 19 (yield of 79%) was obtained in the same manner as in Synthesis Example 1, except that Intermediate A was used instead of Intermediate A-1.

[0320] ¹H NMR (300 MHz, CDCl₃) 8.49 (2H), 8.32 (3H), 8.15 (1H), 8.01-7.95 (3H), 7.86-7.81 (5H), 7.79 (2H), 7.64-7.56 (4H), 7.53-7.42 (7H), 7.39 (1H), 7.24-7.19 (2H), 6.95 (1H)

Synthesis Example 4

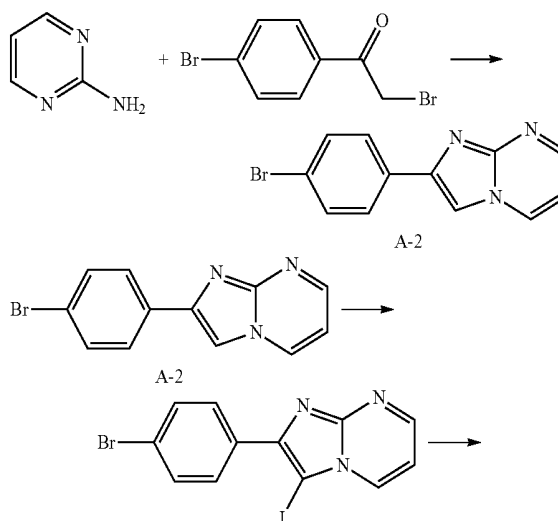
Synthesis of Compound 18

[0321] Compound 18 (yield of 83%) was obtained in the same manner as in Synthesis Example 2, except that Intermediate A was used instead of Intermediate A-1.

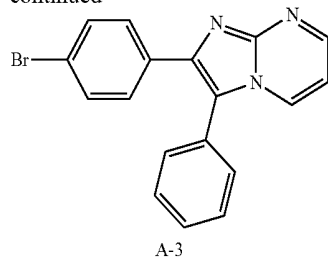
[0322] ¹H NMR (300 MHz, CDCl₃) 8.54-8.53 (2H), 8.42 (1H), 8.15-8.09 (3H), 7.95-7.90 (6H), 7.86-7.79 (2H), 7.69 (2H), 7.63-7.48 (5H), 7.39-7.34 (3H), 7.25-7.17 (5H), 7.03 (1H)

Synthesis Example 5

Synthesis of Compound 7

[0323]

-continued



[0324] Compound 7 (yield of 72%) was obtained in the same manner as in Synthesis Example 1, except that Intermediate A-3 was used instead of Intermediate A-1.

[0325] ^1H NMR (300 MHz, CDCl_3) 8.79-8.75 (2H), 8.53 (1H), 8.30 (3H), 8.16 (1H), 7.98-7.93 (3H), 7.87-7.81 (6H), 7.78-7.63 (6H), 7.59-7.43 (10H), 7.39-7.34 (2H), 7.05 (1H)

Example 1

[0326] An ITO glass substrate (Corning Inc.) with an ITO layer having a thickness of $15\ \Omega/\text{cm}^2$ (500 Å) was cut to a size of 50 mm×50 mm×0.5 mm and sonicated in isopropyl alcohol and water, each for 10 minutes, and then subjected to UV-radiation for 10 minutes and cleaned by exposure to ozone. Then the ITO glass substrate was placed in a vacuum desposition apparatus.

[0327] 2-TNATA was deposited on the ITO anode glass substrate to form a hole injection layer having a thickness of 600 Å, and then NPB was deposited on the hole injection layer to form a hole transport layer having a thickness of 300 Å. After that, Compound 1 (host) and $\text{Ir}(\text{ppy})_3$ (dopant) were co-deposited in a weight ratio of 87:13 on the hole transport layer to form an emission layer having a thickness of 300 Å.

[0328] Alq_3 was deposited on the emission layer to form an electron transport layer having a thickness of 300 Å, and Al was deposited on the electron transport layer to form a cathode having a thickness of 1,200 Å, thereby manufacturing an organic light-emitting device.

Example 2

[0329] An organic light-emitting device was manufactured in the same manner as in Example 1, except that in forming an emission layer, Compound 10 was used instead of Compound 1.

Example 3

[0330] An organic light-emitting device was manufactured in the same manner as in Example 1, except that in forming an emission layer, Compound 19 was used instead of Compound 1.

Example 4

[0331] An organic light-emitting device was manufactured in the same manner as in Example 1, except that in forming an emission layer, Compound 18 was used instead of Compound 1.

Example 5

[0332] An organic light-emitting device was manufactured in the same manner as in Example 1, except that in forming an emission layer, Compound 7 was used instead of Compound 1.

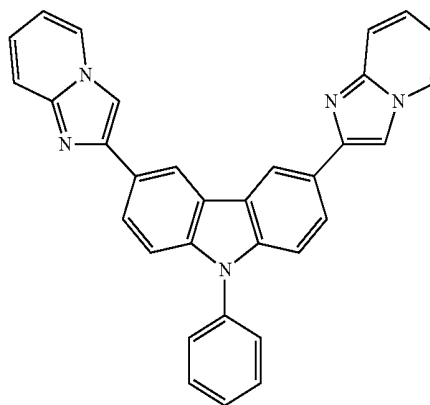
Comparative Example 1

[0333] An organic light-emitting device was manufactured in the same manner as in Example 1, except that in forming an emission layer, as a host, CBP was used instead of Compound 1.

Comparative Example 2

[0334] An organic light-emitting device was manufactured in the same manner as in Example 1, except that in forming an emission layer, as a host, Compound A below was used instead of Compound 1.

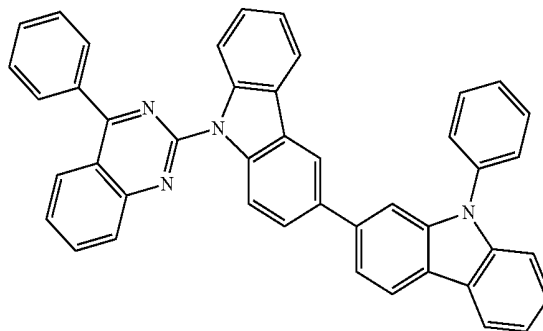
<Compound A>



Comparative Example 3

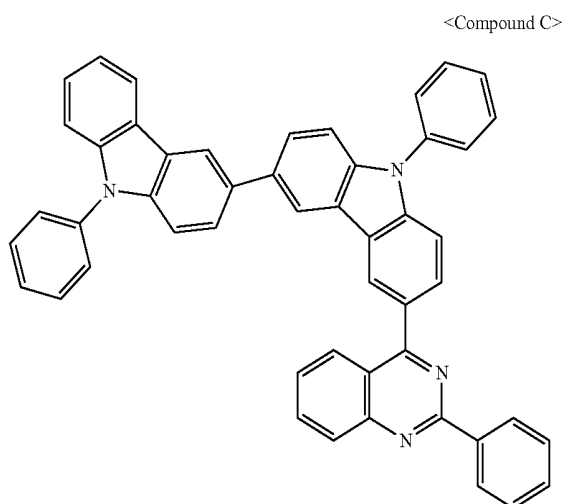
[0335] An organic light-emitting device was manufactured in the same manner as in Example 1, except that in forming an emission layer, as a host, Compound B below was used instead of Compound 1.

<Compound B>



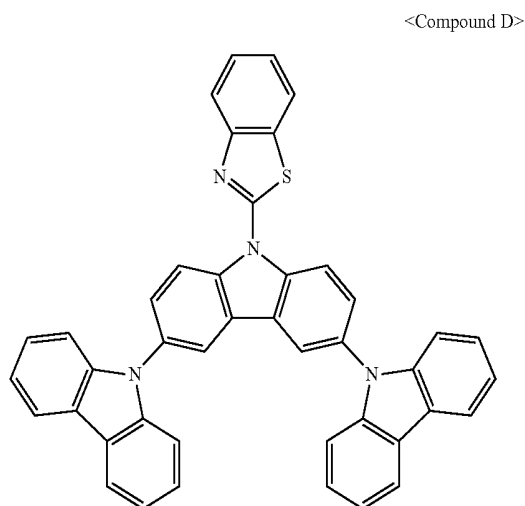
Comparative Example 4

[0336] An organic light-emitting device was manufactured in the same manner as in Example 1, except that in forming an emission layer, as a host, Compound C below was used instead of Compound 1.



Comparative Example 5

[0337] An organic light-emitting device was manufactured in the same manner as in Example 1, except that in forming an emission layer, as a host, Compound D below was used instead of Compound 1.



Evaluation Example 1

[0338] The driving voltage, current density, efficiency, color purity, and T95 lifetime of the organic light-emitting devices manufactured according to Examples 1-5 and Comparative Examples 1-5 were measured by using a Keithley 236 SMU and a brightness photometer PR650, and results thereof are shown in Table 1. T95 lifetime refers to a period of time that is taken until the brightness of the organic light-emitting device was 95% of its initial brightness, at the current density of 10 mA/cm².

TABLE 1

	Emission layer host	Driving voltage (V)	Current Density (mA/cm ²)	Efficiency (cd/A)	T95 lifetime	
					CIE	(hr)
Example 1	Compound 1	3.7	10	60.8	(0.32, 0.62)	289
Example 2	Compound 10	4.2	10	62.3	(0.33, 0.60)	377
Example 3	Compound 19	4.2	10	61.5	(0.31, 0.62)	326
Example 4	Compound 18	4.4	10	68.2	(0.32, 0.63)	399
Example 5	Compound 7	4.0	10	65.5	(0.32, 0.62)	315
Comparative Example 1	CBP	5.3	10	52.3	(0.33, 0.61)	185
Comparative Example 2	Compound A	5.1	10	51.2	(0.32, 0.63)	225
Comparative Example 3	Compound B	5.5	10	46.5	(0.31, 0.62)	264
Comparative Example 4	Compound C	4.9	10	43.1	(0.33, 0.61)	219
Comparative Example 5	Compound D	5.8	10	53.7	(0.32, 0.61)	172

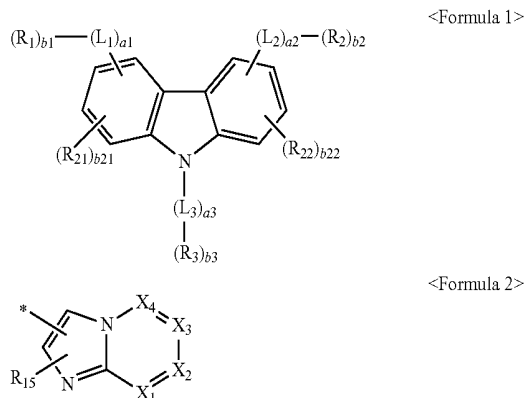
[0339] From Table 1, it may be seen that the driving voltage, brightness, efficiency, and T95 lifetime of the organic light-emitting devices manufactured according to Examples 1-5 were improved, relative to the driving voltage, brightness, efficiency, and T95 lifetime of the organic light-emitting devices manufactured according to Comparative Examples 1-5.

[0340] As described above, according to the one or more of the above exemplary embodiments, an organic light-emitting device including the carbazole-based compound according to an embodiment may have a low driving voltage, high efficiency, high brightness, and long lifespan.

[0341] Example embodiments have been disclosed herein, and although specific terms are employed, they are used and are to be interpreted in a generic and descriptive sense only and not for purpose of limitation. In some instances, as would be apparent to one of ordinary skill in the art as of the filing of the present application, features, characteristics, and/or elements described in connection with a particular embodiment may be used singly or in combination with features, characteristics, and/or elements described in connection with other embodiments unless otherwise specifically indicated. Accordingly, it will be understood by those of skill in the art that various changes in form and details may be made without departing from the spirit and scope of the present invention as set forth in the following claims.

What is claimed is:

1. A carbazole-based compound represented by Formula 1 below:



wherein, in Formulae 1 and 2,

X₁ is C(R₁₁) or N, X₂ is C(R₁₂) or N, X₃ is C(R₁₃) or N, and X₄ is C(R₁₄) or N;

R₁₁ to R₁₅, R₂₁ and R₂₂ are each independently selected from a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a substituted or unsubstituted C₁-C₆₀ alkyl group, a substituted or unsubstituted C₂-C₆₀ alkenyl group, a substituted or unsubstituted C₂-C₆₀ alkynyl group, a substituted or unsubstituted C₁-C₆₀ alkoxy group, a substituted or unsubstituted C₃-C₁₀ cycloalkyl group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkyl group, a substituted or unsubstituted C₃-C₁₀ cycloalkenyl group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkenyl group, a substituted or unsubstituted C₆-C₆₀ aryl group, a substituted or unsubstituted C₆-C₆₀ aryloxy group, a substituted or unsubstituted C₆-C₆₀ arylthio group, a substituted or unsubstituted C₁-C₆₀ heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, —Si(Q₁)(Q₂)(Q₃), —B(Q₄)(Q₅) and —N(Q₆)(Q₇);

b₂₁ and b₂₂ are each independently selected from 0, 1, 2, and 3;

L₁ to L₃ are each independently selected from a substituted or unsubstituted C₃-C₁₀ cycloalkylene group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkylene group, a substituted or unsubstituted C₃-C₁₀ cycloalkenylene group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkenylene group, a substituted or unsubstituted C₆-C₆₀ arylene group, a substituted or unsubstituted C₁-C₆₀ heteroarylene group, a substituted or unsubstituted divalent non-aromatic condensed polycyclic group, and a substituted or unsubstituted divalent non-aromatic condensed heteropolycyclic group;

a₁ to a₃ are each independently selected from 0, 1, 2, and 3, and when a₁ is 2 or more, two or more of L₁ are present and the two or more L₁ are identical or different, and when a₂ is 2 or more, two or more L₂ are present and the

two or more L₂ are identical or different, when a₃ is 2 or more, two or more L₃ are present and two or more L₃ are identical or different;

R₁ is a substituted or unsubstituted carbazolyl group;

R₂ and R₃ are each independently selected from a group represented by Formula 2, a hydrogen, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a substituted or unsubstituted C₁-C₆₀ alkyl group, a substituted or unsubstituted C₂-C₆₀ alkenyl group, a substituted or unsubstituted C₂-C₆₀ alkynyl group, a substituted or unsubstituted C₁-C₆₀ alkoxy group, a substituted or unsubstituted C₃-C₁₀ cycloalkyl group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkyl group, a substituted or unsubstituted C₃-C₁₀ cycloalkenyl group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkenyl group, a substituted or unsubstituted C₆-C₆₀ aryl group, a substituted or unsubstituted C₁-C₆₀ heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, —Si(Q₁)(Q₂)(Q₃), —B(Q₄)(Q₅), and —N(Q₆)(Q₇), wherein the substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group does not include a group represented by Formula 2;

at least one of R₂ and R₃ is a group represented by Formula 2;

b₁ to b₃ are each independently an integer of 1 to 3;

at least one substituent of the substituted carbazolyl group, substituted C₃-C₁₀ cycloalkylene group, substituted C₁-C₁₀ heterocycloalkylene group, substituted C₃-C₁₀ cycloalkenylene group, substituted C₁-C₁₀ heterocycloalkenylene group, substituted C₆-C₆₀ arylene group, substituted C₁-C₆₀ heteroarylene group, substituted divalent non-aromatic condensed polycyclic group, substituted divalent non-aromatic condensed heteropolycyclic group, substituted C₁-C₆₀ alkyl group, substituted C₂-C₆₀ alkenyl group, substituted C₂-C₆₀ alkynyl group, substituted C₁-C₆₀ alkoxy group, substituted C₃-C₁₀ cycloalkyl group, substituted C₁-C₁₀ heterocycloalkyl group, substituted C₃-C₁₀ cycloalkenyl group, substituted C₁-C₁₀ heterocycloalkenyl group, substituted C₆-C₆₀ aryl group, substituted C₆-C₆₀ aryloxy group, substituted C₆-C₆₀ arylthio group, substituted C₁-C₆₀ heteroaryl group, substituted monovalent non-aromatic condensed polycyclic group, and substituted monovalent non-aromatic condensed heteropolycyclic group is selected from

a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group;

a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group and a C₁-C₆₀ alkoxy group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt

thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, —Si(Q₁₁)(Q₁₂)(Q₁₃), —B(Q₁₄)(Q₁₅) and —N(Q₁₆)(Q₁₇);

a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic heterocondensed polycyclic group;

a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group and a monovalent non-aromatic condensed heteropolycyclic group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, —Si(Q₂₁)(Q₂₂)(Q₂₃), —B(Q₂₄)(Q₂₅) and —N(Q₂₆)(Q₂₇); and

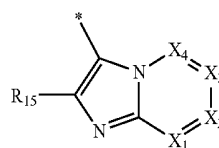
Si(Q₃₁)(Q₃₂)(Q₃₃), —B(Q₃₄)(Q₃₅) and —N(Q₃₆)(Q₃₇);

wherein Q₁ to Q₇, Q₁₁ to Q₁₇, Q₂₁ to Q₂₇, and Q₃₁ to Q₃₇ are each independently selected from a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group, and

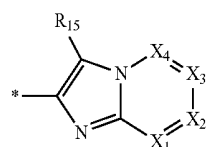
wherein * is a binding site to a neighboring atom.

2. The carbazole-based compound as claimed in claim 1, wherein the group represented by Formula 2 is represented by Formula 2A or 2B below:

<Formula 2A>



<Formula 2B>



in Formulae 2A and 2B, X₁ to X₄ and R₁₅ are the same as those defined with respect to Formula 2, and * is a binding site to a neighboring atom.

3. The carbazole-based compound as claimed in claim 1, wherein X₁ is C(R₁₁), X₂ is C(R₁₂), X₃ is C(R₁₃), and X₄ is C(R₁₄).

4. The carbazole-based compound as claimed in claim 1, wherein X₁ is N, X₂ is C(R₁₂), X₃ is C(R₁₃), and X₄ is C(R₁₄).

5. The carbazole-based compound as claimed in claim 1, wherein R₁₁ to R₁₅, R₂₁, and R₂₂ are each independently selected from:

a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C₁-C₂₀ alkyl group, and a C₁-C₂₀ alkoxy group;

a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothienophenyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzooxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothienophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group; and

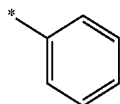
a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzoimidazolyl group, a benzofuranyl group, a benzothienophenyl group, an isobenzothiazolyl group, a ben-

zoxazolyl group, an isobenzoxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group, each substituted with at least one selected from a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, C₁-C₂₀ alkyl group, C₁-C₂₀ alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spirofluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl group, a benzofuranyl group, a benzothiophenyl group, an isobenzothiiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group.

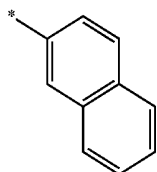
6. The carbazole-based compound as claimed in claim 1, wherein R₁₁ to R₁₅, R₂₁, and R₂₂ are each independently selected from:

a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, C₁-C₂₀ alkyl group and C₁-C₂₀ alkoxy group; and

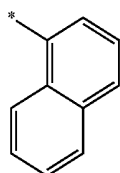
a group represented by one of Formulae 7-1 to 7-45 below:



Formula 7-1

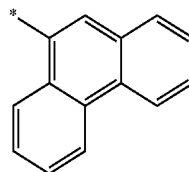


Formula 7-2

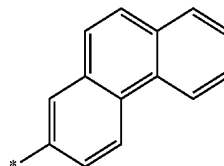


Formula 7-3

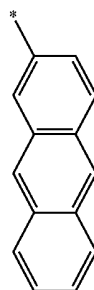
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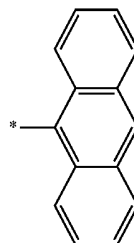
Formula 7-4



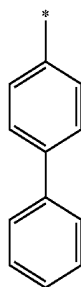
Formula 7-5



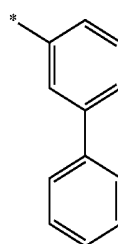
Formula 7-6



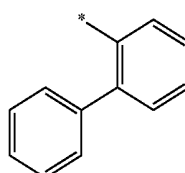
Formula 7-7



Formula 7-8

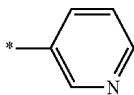
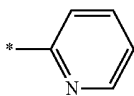
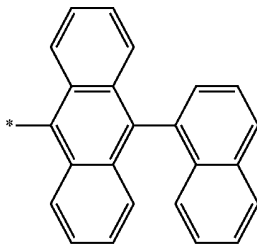
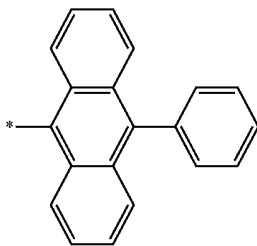
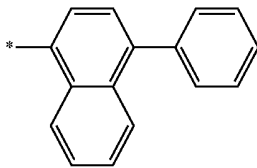
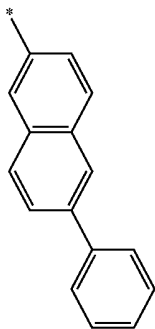
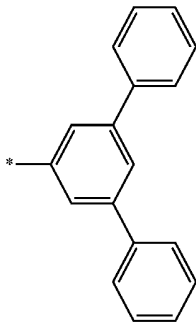


Formula 7-9



Formula 7-10

-continued



Formula 7-11

Formula 7-12

Formula 7-13

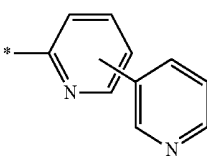
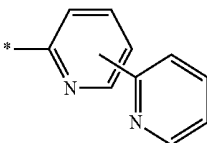
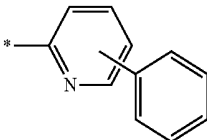
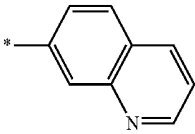
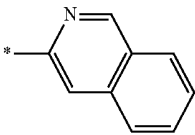
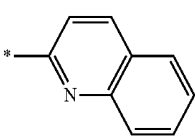
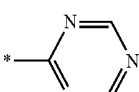
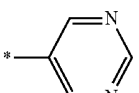
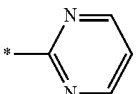
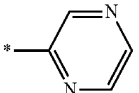
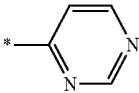
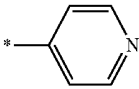
Formula 7-14

Formula 7-15

Formula 7-16

Formula 7-17

-continued



Formula 7-18

Formula 7-19

Formula 7-20

Formula 7-21

Formula 7-22

Formula 7-23

Formula 7-24

Formula 7-25

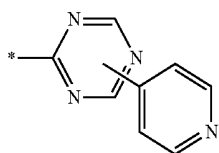
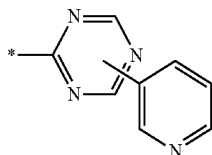
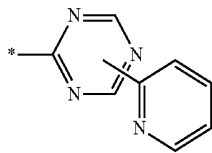
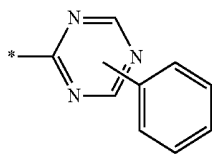
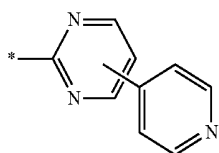
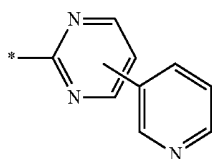
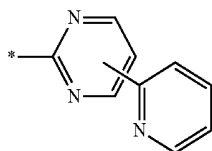
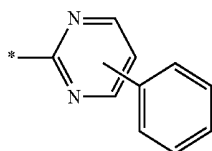
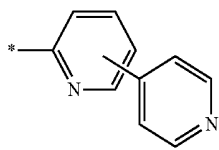
Formula 7-26

Formula 7-27

Formula 7-28

Formula 7-29

-continued



Formula 7-30

Formula 7-31

Formula 7-32

Formula 7-33

Formula 7-34

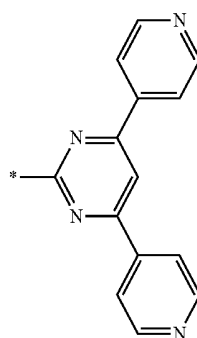
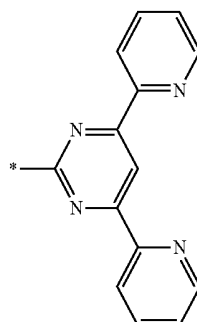
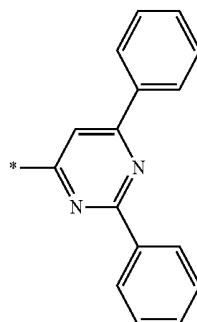
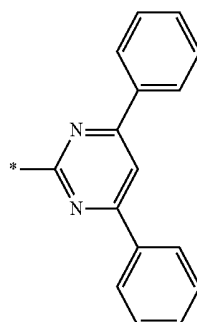
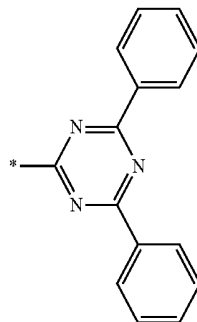
Formula 7-35

Formula 7-36

Formula 7-37

Formula 7-38

-continued



Formula 7-39

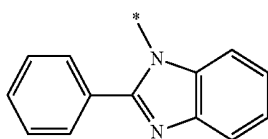
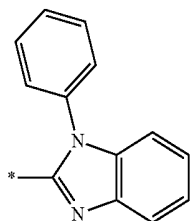
Formula 7-40

Formula 7-41

Formula 7-42

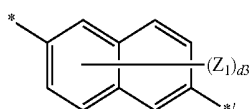
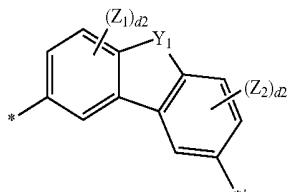
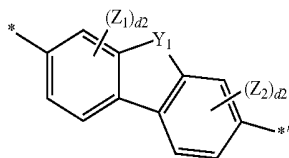
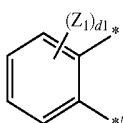
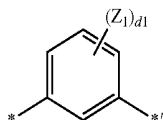
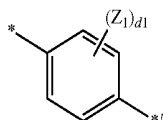
Formula 7-43

-continued



wherein, in Formulae 7-1 to 7-45, * is a binding site to a neighboring atom.

7. The carbazole-based compound as claimed in claim 1, wherein L_1 to L_3 are each independently a group represented by one of Formulae 3-1 to 3-33 below:



Formula 7-44

Formula 7-45

Formula 3-1

Formula 3-2

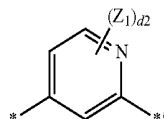
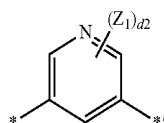
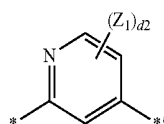
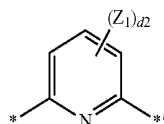
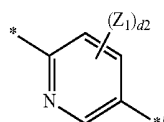
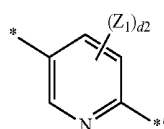
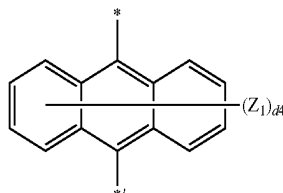
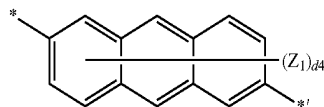
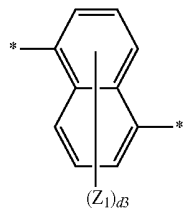
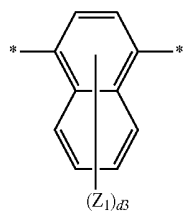
Formula 3-3

Formula 3-4

Formula 3-5

Formula 3-6

-continued



Formula 3-7

Formula 3-8

Formula 3-9

Formula 3-10

Formula 3-11

Formula 3-12

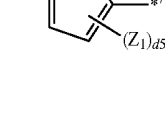
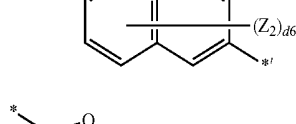
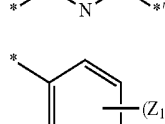
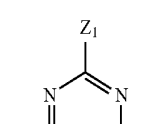
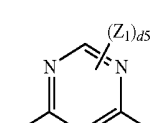
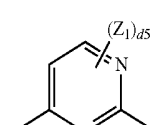
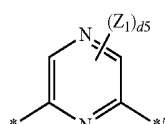
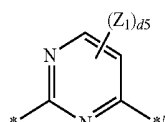
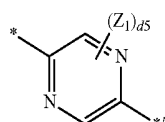
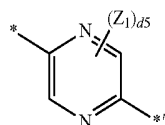
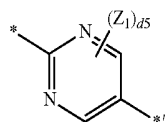
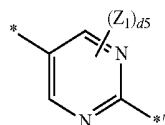
Formula 3-13

Formula 3-14

Formula 3-15

Formula 3-16

-continued



Formula 3-17

Formula 3-18

Formula 3-19

Formula 3-20

Formula 3-21

Formula 3-22

Formula 3-23

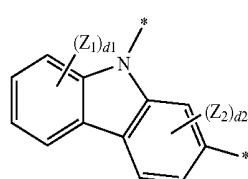
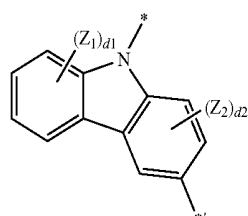
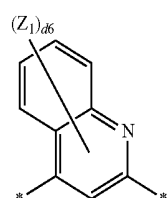
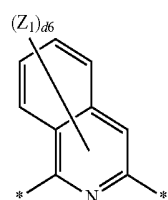
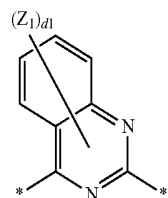
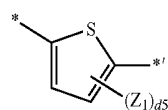
Formula 3-24

Formula 3-25

Formula 3-26

Formula 3-27

-continued



Formula 3-28

Formula 3-29

Formula 3-30

Formula 3-31

Formula 3-32

Formula 3-33

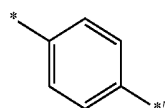
wherein, in Formulae 3-1 to 3-33,

Y_1 is O, S, $C(Z_3)(Z_4)$, $N(Z_5)$, or $Si(Z_6)(Z_7)$;

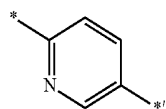
Z_1 to Z_7 are each independently selected from a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 to C_{20} alkyl group, a C_1 to C_{20} alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, and a triazinyl group;

d1 is an integer of 1 to 4, d2 is an integer of 1 to 3, d3 is an integer of 1 to 6, d4 is an integer of 1 to 8, d5 is 1 or 2, and d6 is an integer of 1 to 5, and * and *' indicate binding sites to a neighboring atom.

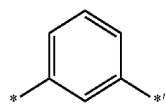
8. The carbazole-based compound as claimed in claim 1, wherein L_1 to L_3 are each independently a group represented by one of Formulae 4-1 to 4-24 below:



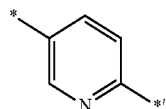
Formula 4-1



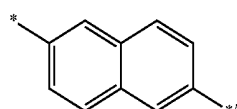
Formula 4-2



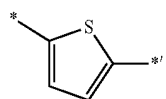
Formula 4-3



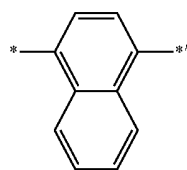
Formula 4-4



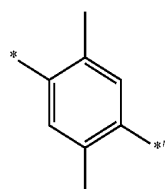
Formula 4-5



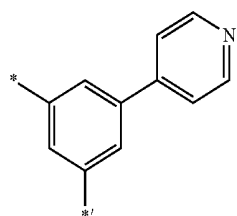
Formula 4-6



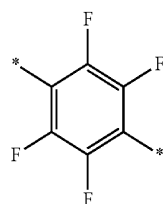
Formula 4-7



Formula 4-8

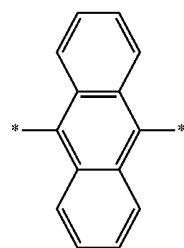


Formula 4-9

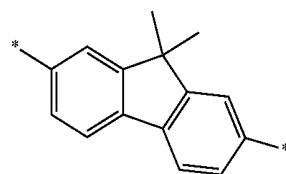


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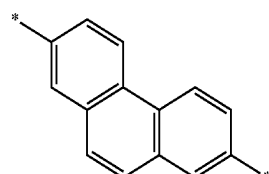
Formula 4-10



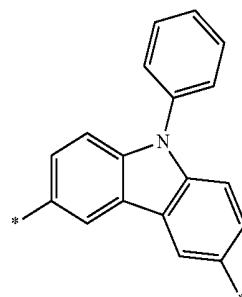
Formula 4-11



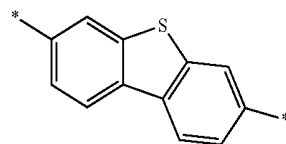
Formula 4-12



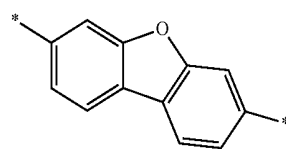
Formula 4-13



Formula 4-14

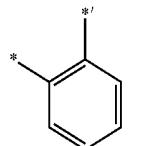
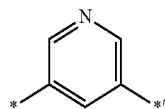
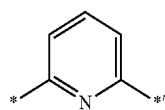
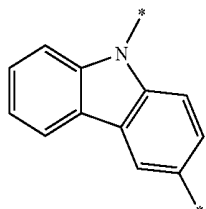
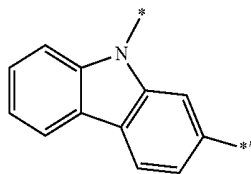
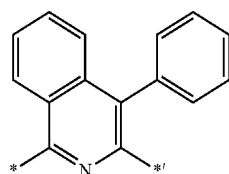
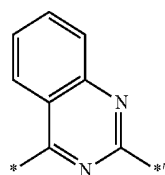
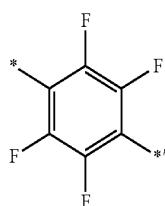


Formula 4-15



Formula 4-16

-continued

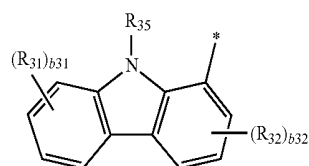


wherein, in Formulae 4-1 to 4-24, * and *' indicate binding sites to a neighboring atom.

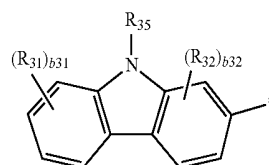
9. The carbazole-based compound as claimed in claim 1, wherein a1 to a3 are each independently 0 or 1.

10. The carbazole-based compound as claimed in claim 1, wherein R₁ is a group represented by one of Formulae 8-1 to 8-5 below:

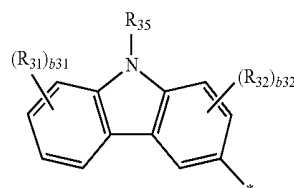
Formula 4-17



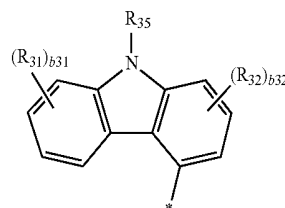
Formula 4-18



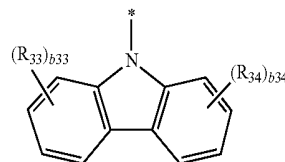
Formula 4-19



Formula 4-20



Formula 4-21



Formula 4-22

wherein in Formulae 8-1 to 8-5,

R₃₁ to R₃₄ are defined the same as R₁₁ of Formula 2;

R₃₅ is selected from a substituted or unsubstituted C₁-C₆₀ alkyl group, a substituted or unsubstituted C₂-C₆₀ alkenyl group, a substituted or unsubstituted C₂-C₆₀ alkynyl group, a substituted or unsubstituted C₁-C₆₀ alkoxy group, a substituted or unsubstituted C₃-C₁₀ cycloalkyl group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkyl group, a substituted or unsubstituted C₃-C₁₀ cycloalkenyl group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkenyl group, a substituted or unsubstituted C₆-C₆₀ aryl group, a substituted or unsubstituted C₁-C₆₀ heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, —Si(Q₁)(Q₂)(Q₃), —B(Q₄)(Q₅), and —N(Q₆)(Q₇);

b31, b33, and b34 are each independently selected from 0, 1, 2, 3, and 4, and b32 is selected from 0, 1, 2, and 3.

11. The carbazole-based compound as claimed in claim 1, wherein:

R₂ is selected from:

a group represented by Formula 2;

a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a

group, an oxadiazolyl group, a triazinyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, and a thiadiazolyl group; and

at least one of R_2 and R_3 is a group represented by Formula 2.

12. The carbazole-based compound as claimed in claim 1, wherein:

R_2 is selected from:

a group represented by Formula 2;

a hydrogen, a deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, and a C_1 - C_{20} alkoxy group; and

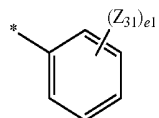
a group represented by one of Formulae 5-1 to 5-26 below;

R_3 is selected from:

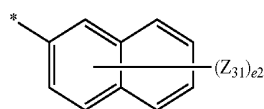
a group represented by Formula 2; and

a group represented by one of Formulae 5-1 to 5-26 below; and

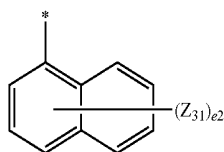
at least one of R_2 and R_3 is a group represented by Formula 2,



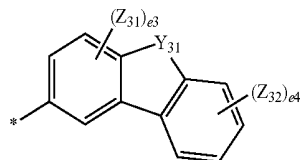
Formula 5-1



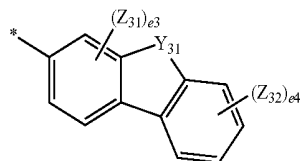
Formula 5-2



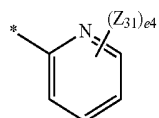
Formula 5-3



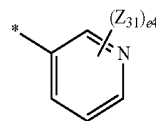
Formula 5-4



Formula 5-5

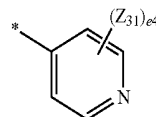


Formula 5-6

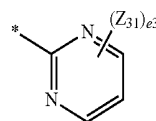


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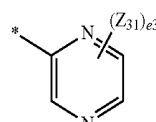
Formula 5-7



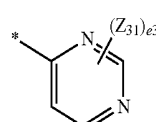
Formula 5-8



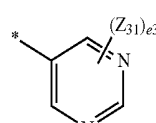
Formula 5-9



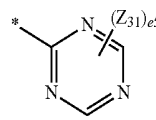
Formula 5-10



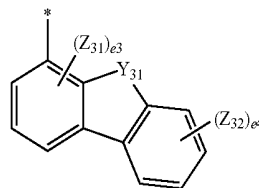
Formula 5-11



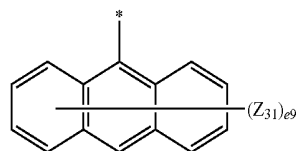
Formula 5-12



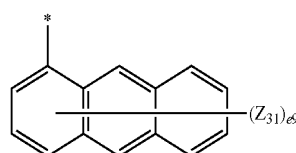
Formula 5-13



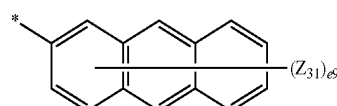
Formula 5-14



Formula 5-15

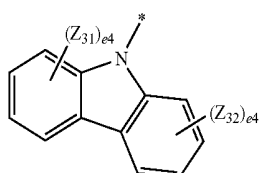
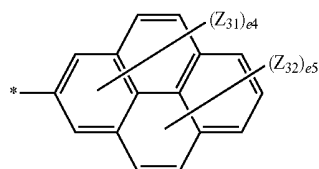
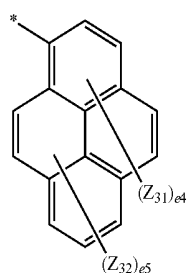
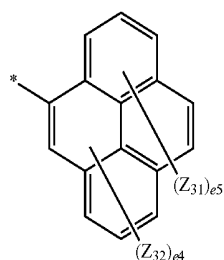
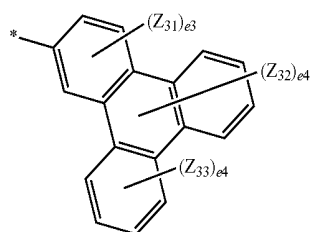
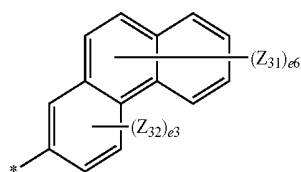
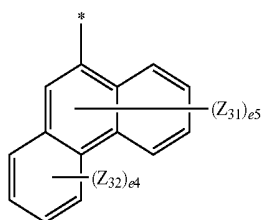


Formula 5-16



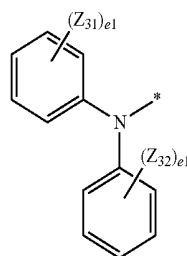
Formula 5-17

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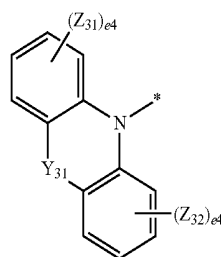
-continued

Formula 5-18



Formula 5-19

Formula 5-20



Formula 5-25

Formula 5-26

wherein, in Formulae 5-1 to 5-26,

Y_{31} is O, S, $C(Z_{34})(Z_{35})$, $N(Z_{36})$, or $Si(Z_{37})(Z_{38})$;

Z_{31} to Z_{38} are each independently selected from:

a hydrogen, a deuterium, $-F$, $-Cl$, $-Br$, $-I$, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, and a C_1 - C_{20} alkoxy group;

a C_1 - C_{20} alkyl group and a C_1 - C_{20} alkoxy group, each substituted with at least one selected from a deuterium, $-F$, $-Cl$, $-Br$, $-I$, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof;

a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a triazinyl group, a benzofuran group, a benzothiophenyl group, a benzocarbazolyl group, and a dibenzocarbazolyl group;

a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a triazinyl group, a benzofuran group, a benzothiophenyl group, a benzocarbazolyl group, and a dibenzocarbazolyl group, each substituted with at least one selected from a deuterium, $-F$, $-Cl$, $-Br$, $-I$, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric

acid or a salt thereof, a C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a phenyl group, and a naphthyl group; and $-\text{Si}(\text{Q}_{31})(\text{Q}_{32})(\text{Q}_{33})$,

wherein Q_{31} to Q_{33} are each independently selected from a hydrogen, a deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, and a triazinyl group;

e_1 is an integer of 1 to 5, e_2 is an integer of 1 to 7, e_3 is an integer of 1 to 3, e_4 is an integer of 1 to 4, e_5 is 1 or 2, e_6 is an integer of 1 to 6, e_9 is an integer of 1 to 9, and * indicates a binding site to a neighboring atom.

13. The carbazole-based compound as claimed in claim 1, wherein:

R_2 is selected from:

a group represented by Formula 2;

a hydrogen, a deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, C_1 - C_{20} alkyl group, and C_1 - C_{20} alkoxy group; and

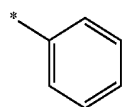
a group represented by one of Formulae 6-1 to 6-35 below;

R_3 is selected from:

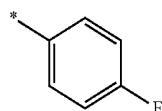
a group represented by Formula 2; and

a group represented by one of Formulae 6-1 to 6-35 below; and

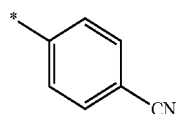
at least one of R_2 and R_3 is a group represented by Formula 2,



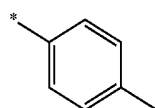
Formula 6-1



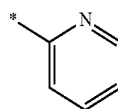
Formula 6-2



Formula 6-3

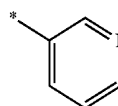


Formula 6-4

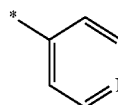


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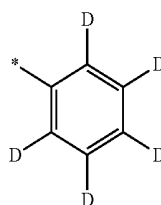
Formula 6-5



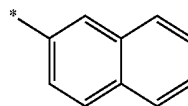
Formula 6-6



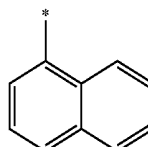
Formula 6-7



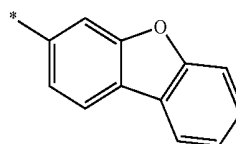
Formula 6-8



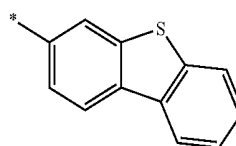
Formula 6-9



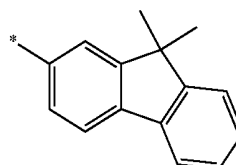
Formula 6-10



Formula 6-11

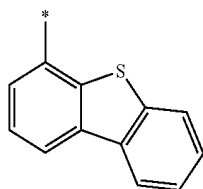
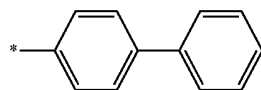
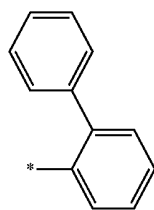
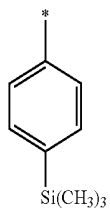
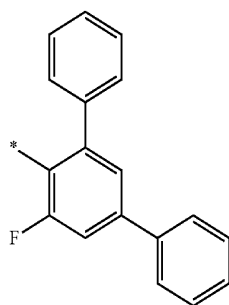
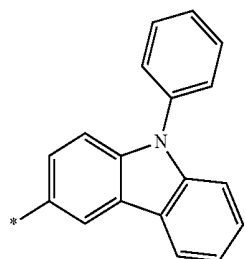


Formula 6-12



Formula 6-13

-continued



Formula 6-14

Formula 6-15

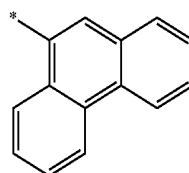
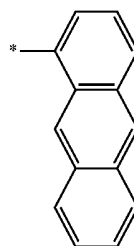
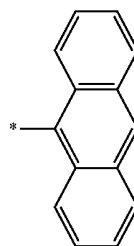
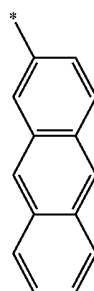
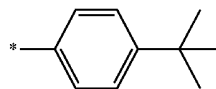
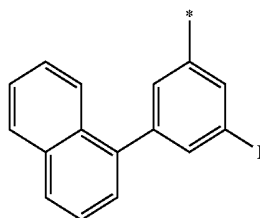
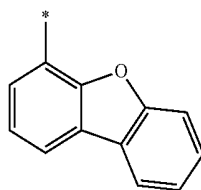
Formula 6-16

Formula 6-17

Formula 6-18

Formula 6-19

-continued



Formula 6-20

Formula 6-21

Formula 6-22

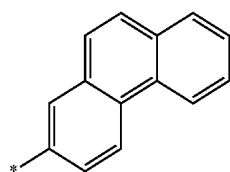
Formula 6-23

Formula 6-24

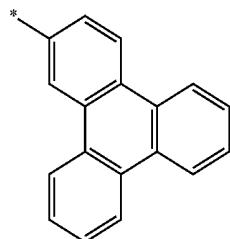
Formula 6-25

Formula 6-26

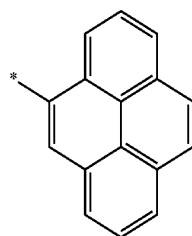
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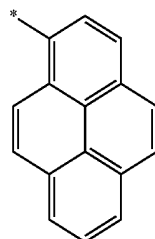
Formula 6-27



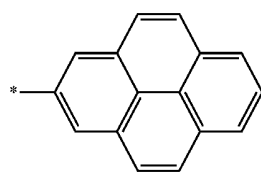
Formula 6-28



Formula 6-29

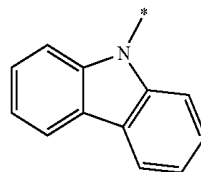


Formula 6-30

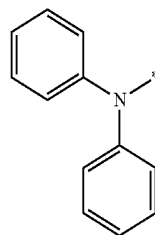


Formula 6-31

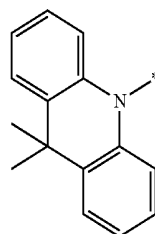
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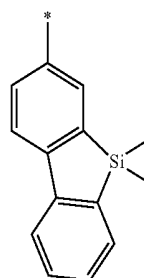
Formula 6-32



Formula 6-33



Formula 6-34

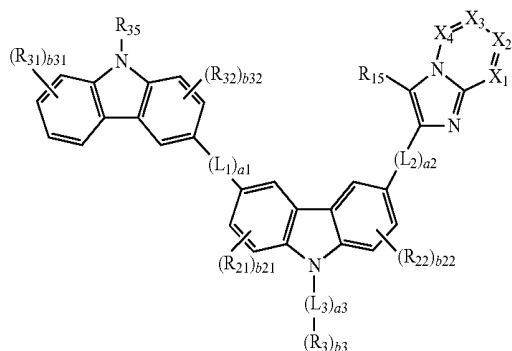


Formula 6-35

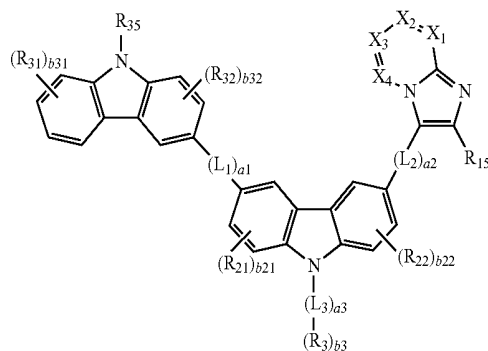
wherein * in Formulae 6-1 to 6-35 indicates a binding site to a neighboring atom.

14. The carbazole-based compound as claimed in claim 1, wherein the carbazole-based compound represented by Formula 1 is represented by one selected from Formulae 1A(1) to 1A(6) and Formulae 1B(1) to 1B(6), below:

<Formula 1A(1)>



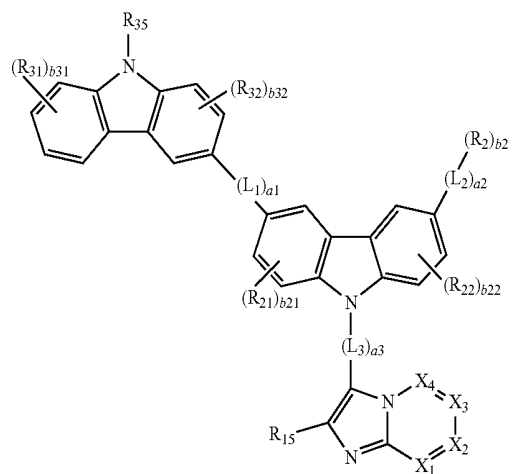
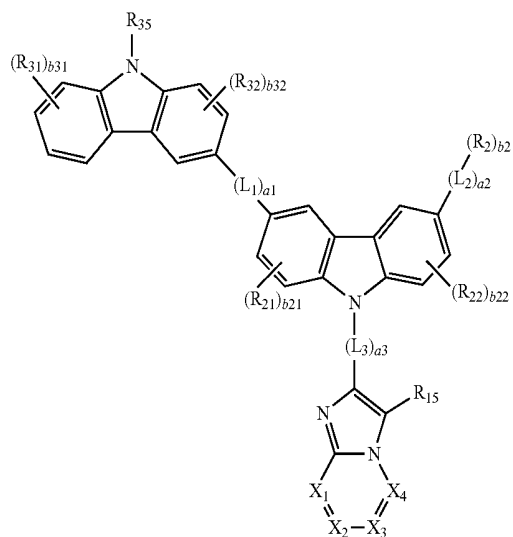
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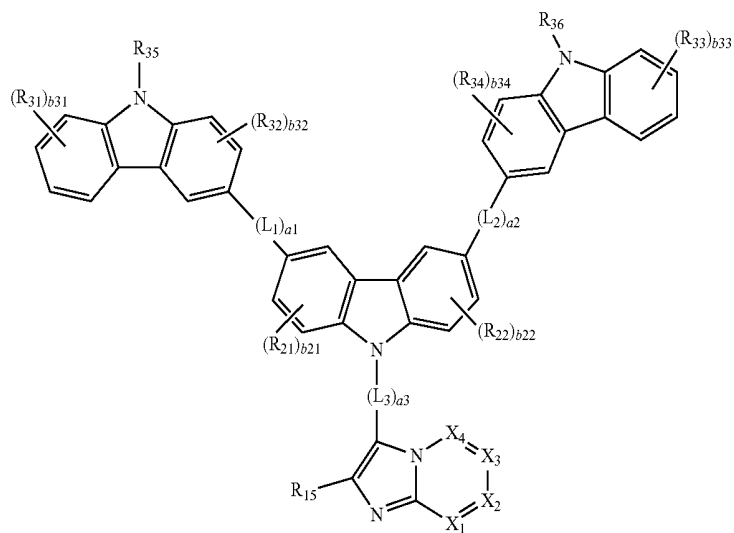
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<Formula 1A(3)>

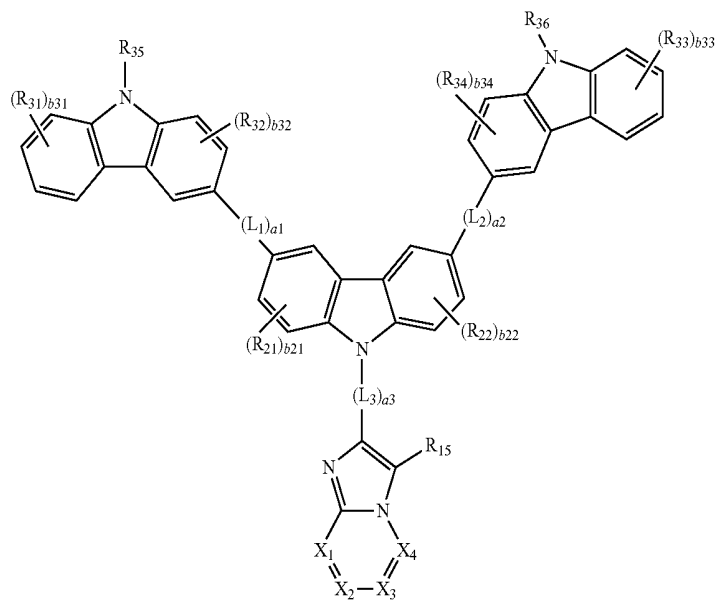
<Formula 1A(4)>



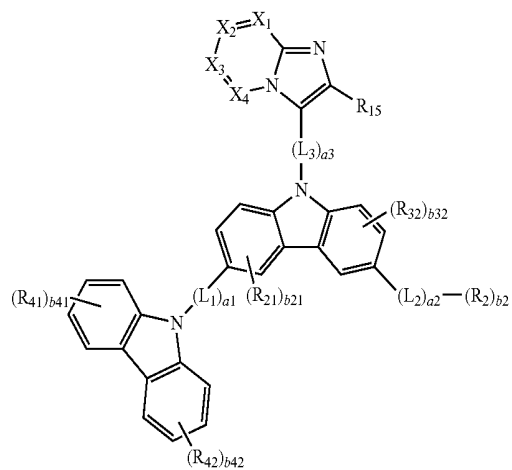
<Formula 1A(5)>



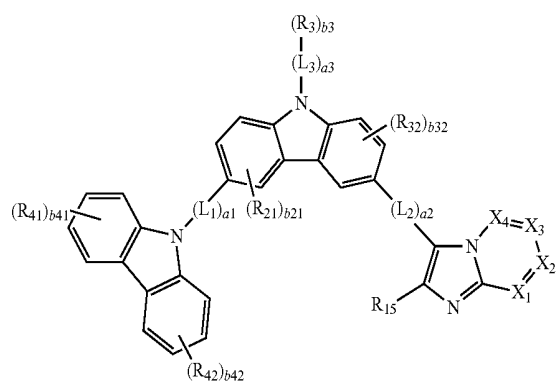
<Formula 1A(6)>



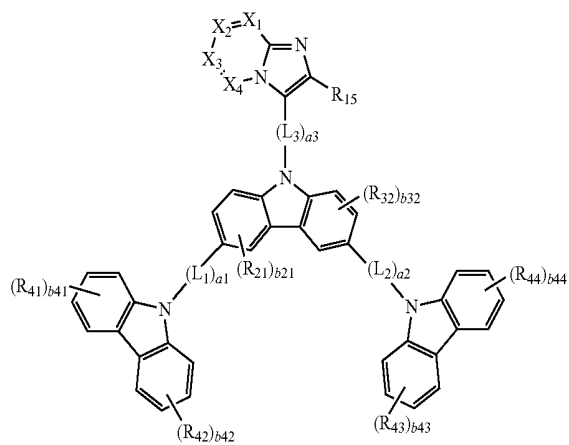
<Formula IB(2)>



<Formula 1B(4)>

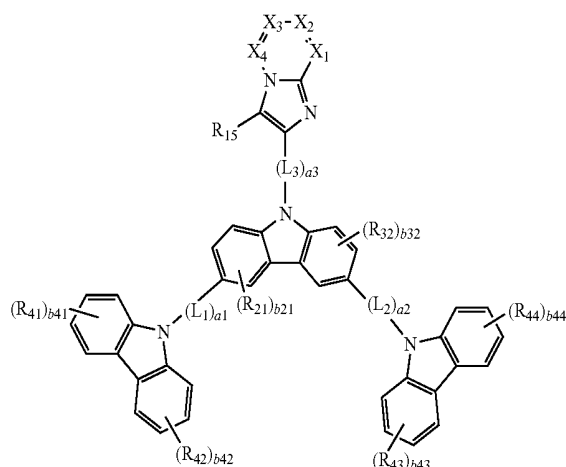


<Formula 1 B(5)>



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<Formula 1B(6)>



wherein, in Formulae 1A(1) to 1A(6) and Formulae 1B(1) to 1B(6),

X_1 to X_4 , R_{15} , R_{21} , R_{22} , b_{21} , b_{22} , L_1 to L_3 , a_1 to a_3 , b_2 , and b_3 are defined the same as those of Formula 1 and Formula 2;

R_2 and R_3 are each independently selected from a group represented by Formula 2, a hydrogen, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxylic acid or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a substituted or unsubstituted C_1 - C_{60} alkyl group, a substituted or unsubstituted C_2 - C_{60} alkenyl group, a substituted or unsubstituted C_2 - C_{60} alkynyl group, a substituted or unsubstituted C_1 - C_{60} alkoxy group, a substituted or unsubstituted C_3 - C_{10} cycloalkyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkyl group, a substituted or unsubstituted C_3 - C_{10} cycloalkenyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenyl group, a substituted or unsubstituted C_6 - C_{60} aryl group, a substituted or unsubstituted C_1 - C_{60} heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, $-\text{Si}(Q_1)(Q_2)(Q_3)$, $-\text{B}(Q_4)(Q_5)$, and $-\text{N}(Q_6)(Q_7)$, wherein the substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group does not include a carbazolyl group and the group represented by Formula 2 above;

R_{31} to R_{34} and R_{41} to R_{44} are defined the same as R_{11} of Formula 2;

R_{35} and R_{36} are each independently selected from a substituted or unsubstituted C_1 - C_{60} alkyl group, a substituted or unsubstituted C_2 - C_{60} alkenyl group, a substituted or unsubstituted C_2 - C_{60} alkynyl group, a substituted or unsubstituted C_1 - C_{60} alkoxy group, a substituted or unsubstituted C_3 - C_{10} cycloalkyl group, a sub-

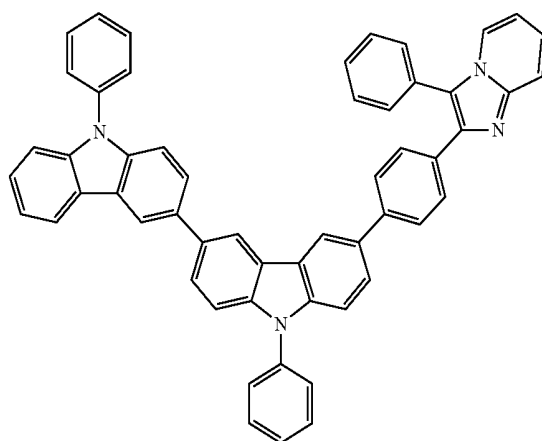
stituted or unsubstituted C_1 - C_{10} heterocycloalkyl group, a substituted or unsubstituted C_3 - C_{10} cycloalkenyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenyl group, a substituted or unsubstituted C_6 - C_{60} aryl group, a substituted or unsubstituted C_1 - C_{60} heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, $-\text{Si}(Q_1)(Q_2)(Q_3)$, $-\text{B}(Q_4)(Q_5)$, and $-\text{N}(Q_6)(Q_7)$;

b_{31} , b_{33} , and b_{41} to b_{44} are each independently selected from 0, 1, 2, 3 and 4, and b_{32} and b_{34} are each independently selected from 0, 1, 2, and 3.

15. The carbazole-based compound as claimed in claim 1, wherein:

X_1 to X_4 are CH; or
 X_1 is N, and X_2 to X_4 are CH.

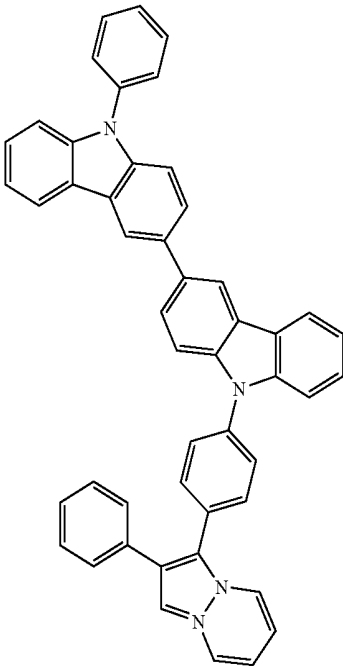
16. The carbazole-based compound as claimed in claim 1, wherein the carbazole-based compound represented by Formula 1 is one of Compounds 1 to 20 below:



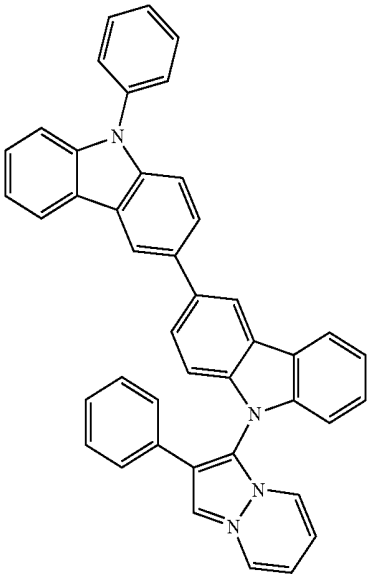
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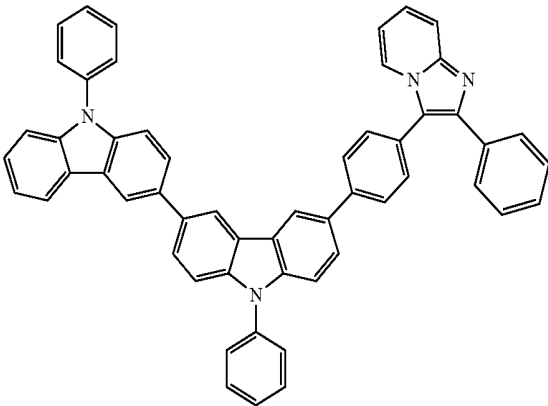
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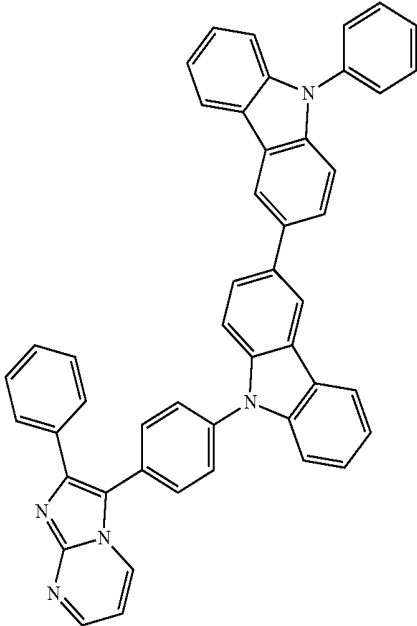


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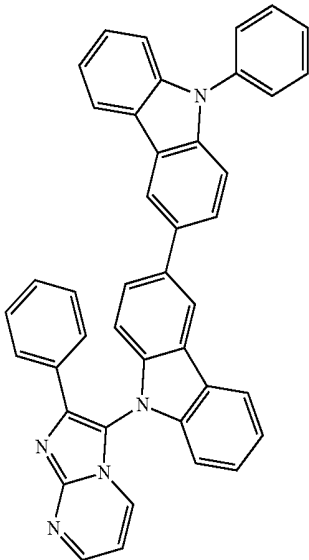


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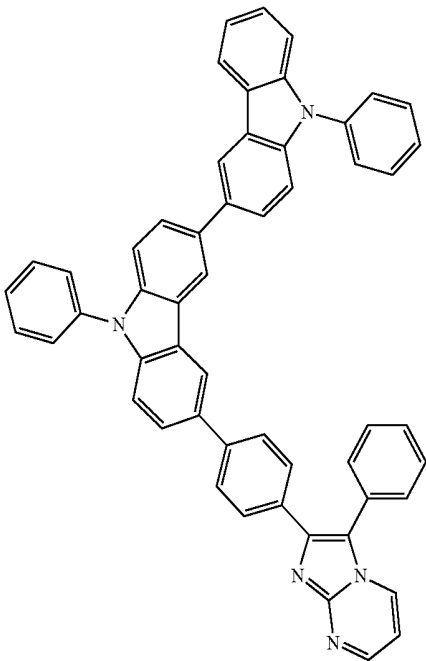
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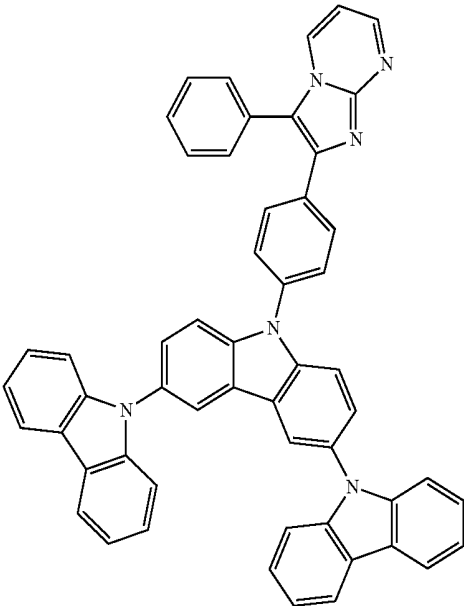


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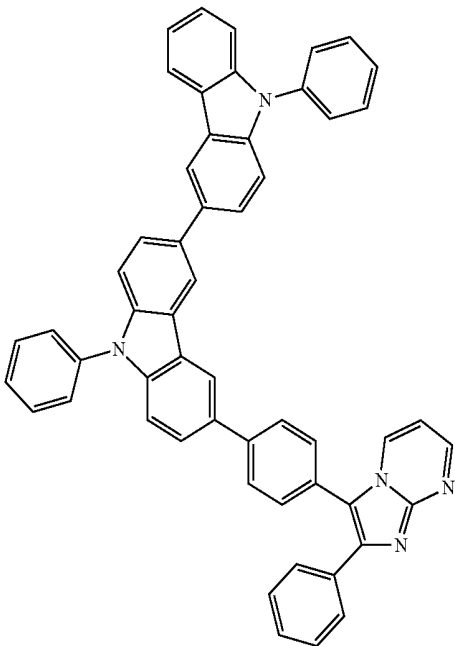


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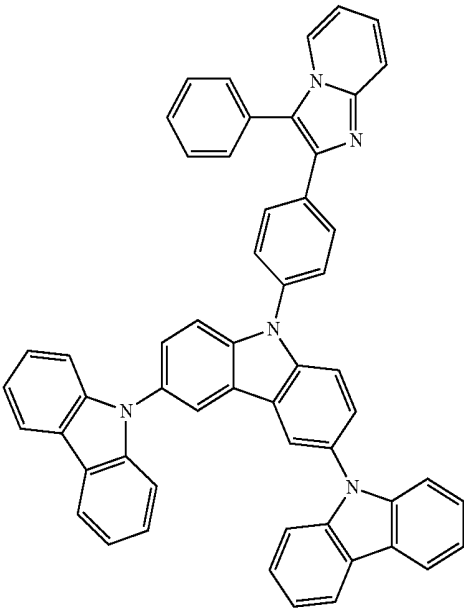
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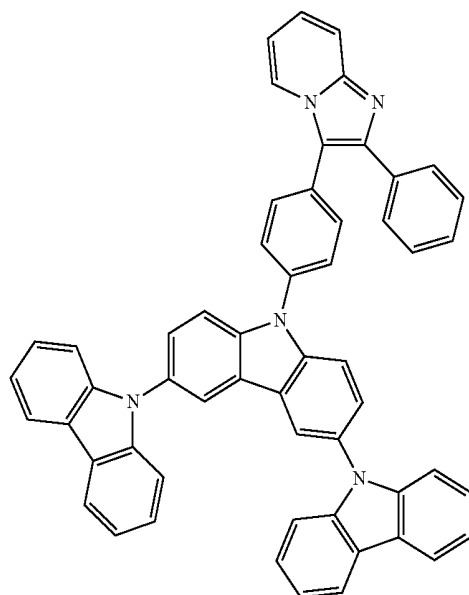


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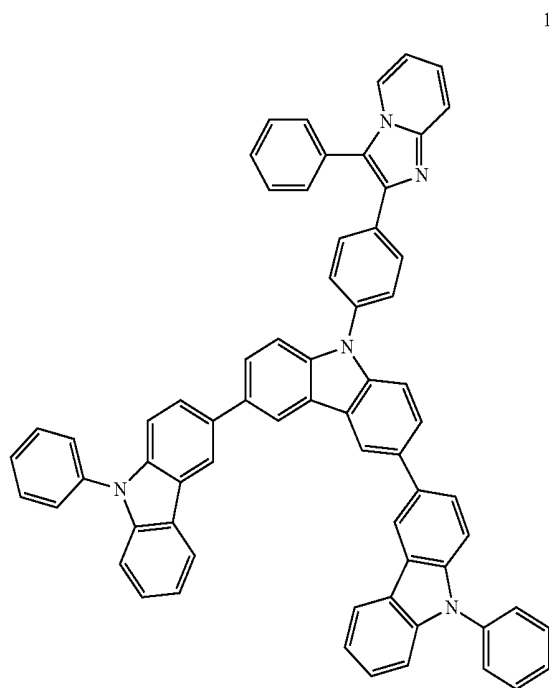


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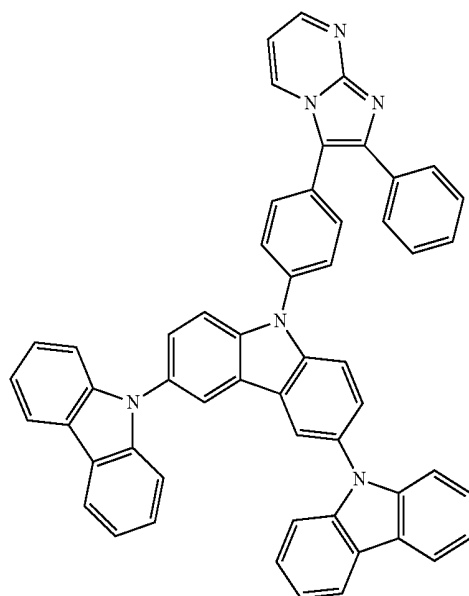
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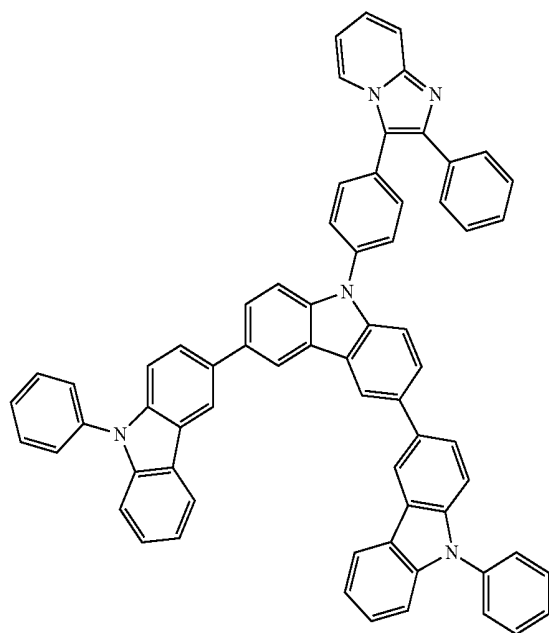
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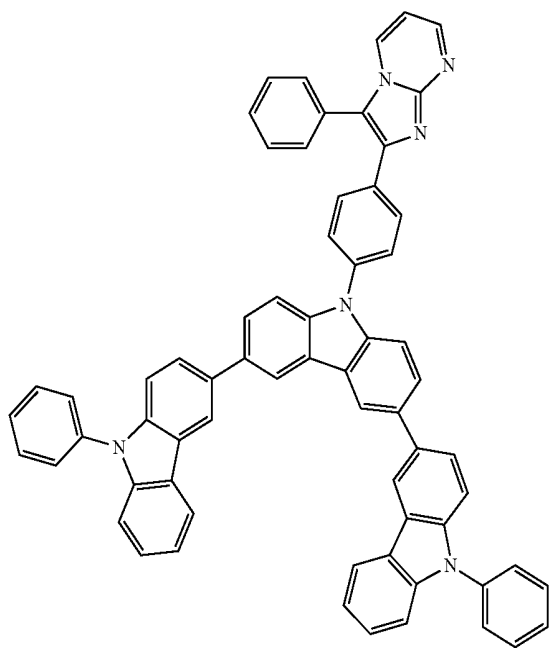


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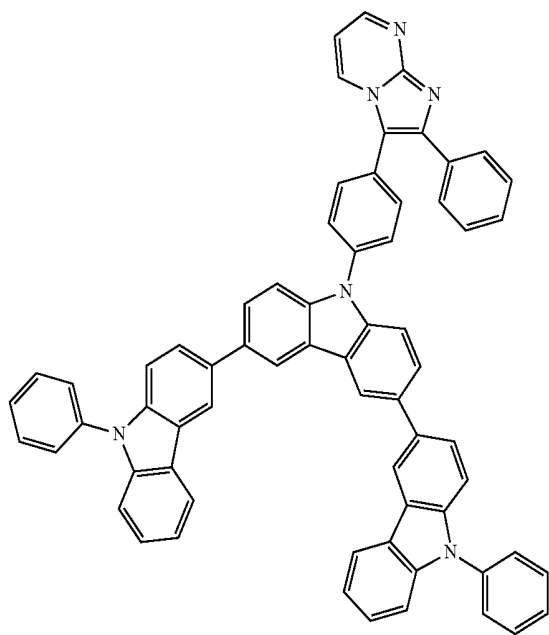
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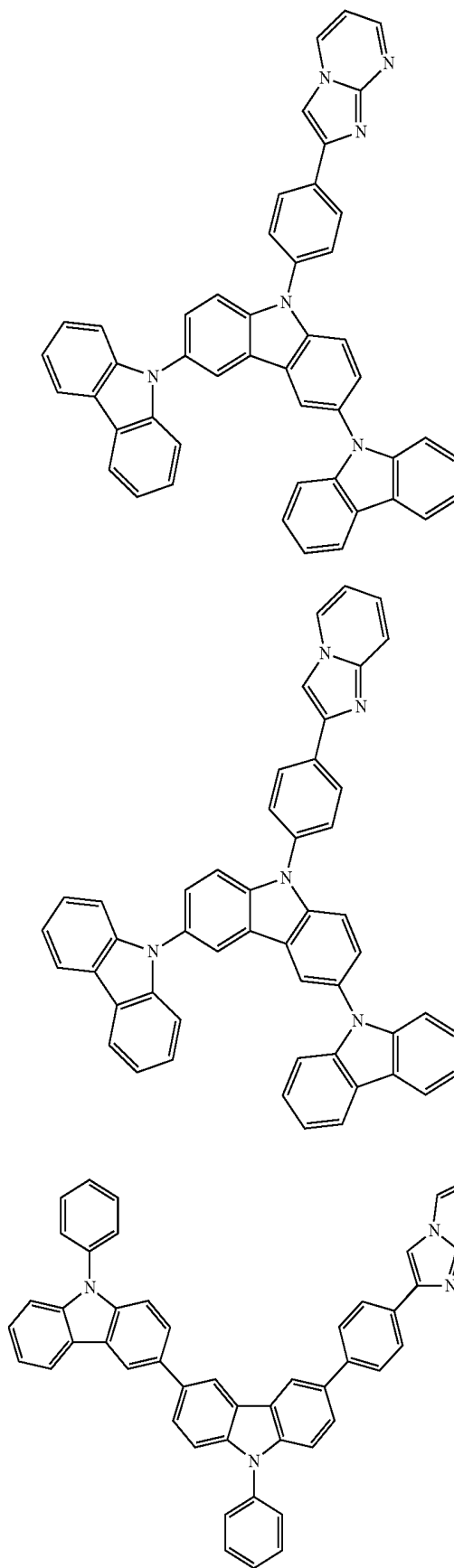


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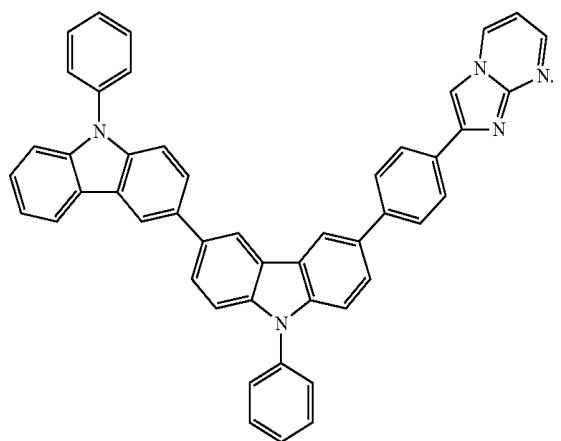
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17. An organic light-emitting device, comprising:
 a first electrode;
 a second electrode facing the first electrode; and
 an organic layer between the first electrode and the second
 electrode and including an emission layer,

wherein the organic layer includes at least one carbazole-
 based compound as claimed in claim 1.

18. The organic light-emitting device as claimed in claim
17, wherein:

the first electrode is an anode,
 the second electrode is a cathode, and
 the organic layer includes:

a hole transport region between the first electrode and
 the emission layer, the hole transport region including
 at least one selected from a hole injection layer, a hole
 transport layer, a buffer layer and an electron blocking
 layer, and

an electron transport region between the emission layer
 and the second electrode, the electron transport region
 including at least one selected from a hole blocking
 layer, an electron transport layer and an electron
 injection layer.

19. The organic light-emitting device as claimed in claim
17, wherein the emission layer includes the at least one car-
 bazole-based compound.

20. The organic light-emitting device as claimed in claim
19, wherein the emission layer further includes a phospho-
 rescent dopant.

* * * * *

专利名称(译)	唑啉类化合物和包括其的有机发光装置		
公开(公告)号	US20160099424A1	公开(公告)日	2016-04-07
申请号	US14/676610	申请日	2015-04-01
[标]申请(专利权)人(译)	三星显示有限公司		
申请(专利权)人(译)	三星DISPLAY CO. , LTD.		
当前申请(专利权)人(译)	三星DISPLAY CO. , LTD.		
[标]发明人	KIM HEEYEON YANG SEUNGGAKE		
发明人	KIM, HEEYEON YANG, SEUNGGAKE		
IPC分类号	H01L51/00 C07D487/04 H01L51/50 C07D471/04		
CPC分类号	H01L51/0072 C07D471/04 H01L51/5012 H01L51/5016 C07D487/04 H01L51/0085		
优先权	1020140132491 2014-10-01 KR		
其他公开文献	US9780313		
外部链接	Espacenet USPTO		

摘要(译)

唑啉类化合物和有机发光装置，唑啉类化合物由化学式1表示

