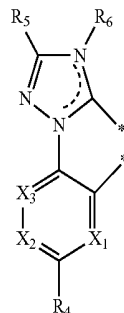


(19) **United States**(12) **Patent Application Publication**
Ko et al.(10) **Pub. No.: US 2018/0114927 A1**(43) **Pub. Date: Apr. 26, 2018**(54) **ORGANOMETALLIC COMPOUND AND ORGANIC LIGHT-EMITTING DEVICE INCLUDING THE SAME** $M(L_1)_{n1}(L_2)_{n2}$

<Formula 1>

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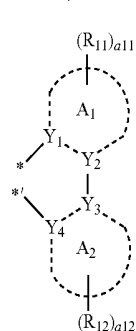
<Formula 2A>

(21) Appl. No.: **15/789,749**(22) Filed: **Oct. 20, 2017**(30) **Foreign Application Priority Data**

Oct. 20, 2016 (KR) 10-2016-0136575

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H01L 51/00 (2006.01)(52) **U.S. Cl.**
CPC **H01L 51/0085** (2013.01); **C07D 213/61** (2013.01); **H01L 51/0067** (2013.01)(57) **ABSTRACT**

An organometallic compound and an organic light-emitting device including the same are provided. The organometallic compound has Formula 1:

wherein M is a third row transition metal, L_1 is a ligand represented by Formula 2A and n_1 is 1, 2, or 3, L_2 is a ligand represented by Formula 2B and n_2 is 0, 1, or 2. When n_1 is two or more, two or more $L_1(s)$ are identical to or different from each other. When n_2 is two or more, two or more $L_2(s)$ are identical to or different from each other. Details about Formula 2A and Formula 2B are provided.

<Formula 2B>

10

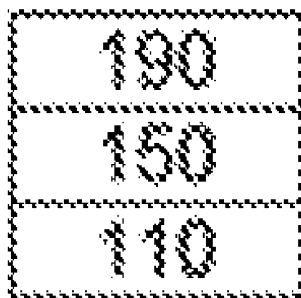
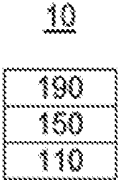


FIG. 1



**ORGANOMETALLIC COMPOUND AND
ORGANIC LIGHT-EMITTING DEVICE
INCLUDING THE SAME**

CROSS-REFERENCE TO RELATED
APPLICATION

[0001] This application claims the benefit of Korean Patent Application No. 10-2016-0136575 filed on Oct. 20, 2016 in the Korean Intellectual Property Office, the disclosure of which is incorporated herein in its entirety by reference.

BACKGROUND

1. Field

[0002] One or more embodiments relate to an organometallic compound and an organic light-emitting device including the same

2. Description of the Related Art

[0003] Organic light-emitting devices are self-emission devices that produce full-color images, and also have wide viewing angles, high contrast ratios, short response times, and excellent characteristics in terms of luminance, driving voltage, and response speed, compared to devices in the art.

[0004] An example of such organic light-emitting devices 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, recombine in the emission layer to produce excitons.

[0005] These excitons transit from an excited state to a ground state, thereby generating light.

SUMMARY

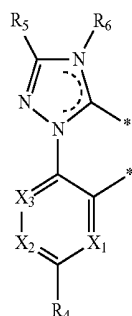
[0006] One or more embodiments include a novel organometallic compound and an organic light-emitting device including the same.

[0007] 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.

[0008] According to one or more embodiments, an organometallic compound is represented by Formula 1:

$M(L_1)_{n1}(L_2)_{n2}$

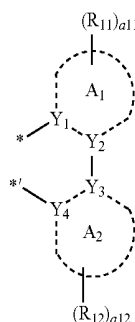
<Formula 1>



<Formula 2A>

-continued

<Formula 2B>



[0009] M in Formula 1 may be a third row transition metal,

[0010] L₁ in Formula 1 may be a ligand represented by Formula 2A, and n₁ may be 1, 2, or 3, wherein, when n₁ is two or more, two or more L₁(s) may be identical to or different from each other,

[0011] L₂ in Formula 1 is a ligand represented by Formula 2B, and n₂ may be 0, 1, or 2, wherein, when n₂ is two or more, two or more L₂(s) may be identical to or different from each other,

[0012] The sum of n₁ and n₂ in Formula 1 may be two or three,

[0013] * and *' in Formulae 2A and 2B each indicate a binding site to M in Formula 1,

[0014] in Formula 2A, X₁ may be N or C(R₁), X₂ may be N or C(R₂), and X₃ may be N or C(R₃),

[0015] in Formula 2B, Y₁ may be carbon, Y₂ and Y₃ may each independently be carbon or nitrogen, Y₄ may be nitrogen, a bond between Y₁ and Y₂ may be a single bond or a double bond, a bond between Y₂ and Y₃ may be a single bond, and a bond between Y₃ and Y₄ may be a single bond or a double bond,

[0016] ring A₁ and ring A₂ in Formula 2B may each be a C₂-C₆₀ heterocyclic group including at least one N as a ring-forming atom,

[0017] R₁ to R₃, R₅, R₆, R₁₁, and R₁₂ in Formulae 2A and 2B may each independently be selected from hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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), —N(Q₁)(Q₂), —B(Q₁)(Q₂), —C(=O)(Q₁), —S(=O)₂(Q₁), and —P(=O)(Q₁)(Q₂).

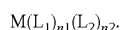
- [0018]** R_4 may be an electron withdrawing group,
- [0019]** R_2 and R_3 in Formula 2A may optionally be linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group,
- [0020]** R_5 and R_6 in Formula 2A may optionally be linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group,
- [0021]** two neighboring groups of a plurality of $R_{11}(s)$ in Formula 2B may optionally be linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group,
- [0022]** two neighboring groups of a plurality of $R_{12}(s)$ in Formula 2B may optionally be linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group,
- [0023]** neighboring R_{11} and R_{12} in Formula 2B may optionally be linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group,
- [0024]** a_{11} and a_{12} in Formula 2B may each independently be an integer from 0 to 10, and
- [0025]** at least one substituent of the substituted C_1 - C_{60} alkyl group, the substituted C_2 - C_{60} alkenyl group, the substituted C_2 - C_{60} alkynyl group, the substituted C_1 - C_{60} alkoxy group, the substituted C_3 - C_{10} cycloalkyl group, the substituted C_1 - C_{10} heterocycloalkyl group, the substituted C_3 - C_{10} cycloalkenyl group, the substituted C_1 - C_{10} heterocycloalkenyl group, the substituted C_6 - C_{60} aryl group, the substituted C_6 - C_{60} aryloxy group, the substituted C_6 - C_{60} arylthio group, the substituted C_1 - C_{60} heteroaryl group, the substituted monovalent non-aromatic condensed polycyclic group, the substituted monovalent non-aromatic condensed heteropolycyclic group, the substituted C_4 - C_{60} carbocyclic group, and the substituted C_2 - C_{60} heterocyclic group may be selected from:
- [0026]** deuterium ($-D$), $-F$, $-Cl$, $-Br$, $-I$, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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;
- [0027]** 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 deuterium, $-F$, $-Cl$, $-Br$, $-I$, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C_3 - C_{10} cycloalkyl group, a C_1 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_1 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, a C_1 - C_{60} heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, $-Si(Q_{11})(Q_{12})(Q_{13})$, $-N(Q_{11})(Q_{12})$, $-B(Q_{11})(Q_{12})$, $-C(=O)(Q_{11})$, $-S(=O)_2(Q_{11})$, and $-P(=O)(Q_{11})(Q_{12})$;
- [0028]** a C_3 - C_{10} cycloalkyl group, a C_1 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_1 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, a C_1 - C_{60} heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group;
- [0029]** a C_3 - C_{10} cycloalkyl group, a C_1 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_1 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, a C_1 - C_{60} 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 deuterium, $-F$, $-Cl$, $-Br$, $-I$, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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 C_3 - C_{10} cycloalkyl group, a C_1 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_1 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, a C_1 - C_{60} heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, $-Si(Q_{21})(Q_{22})(Q_{23})$, $-N(Q_{21})(Q_{22})$, $-B(Q_{21})(Q_{22})$, $-C(=O)(Q_{21})$, $-S(=O)_2(Q_{21})$, and $-P(=O)(Q_{21})(Q_{22})$; and
- [0030]** $Si(Q_{31})(Q_{32})(Q_{33})$, $-N(Q_{31})(Q_{32})$, $-B(Q_{31})(Q_{32})$, $-C(=O)(Q_{31})$, $-S(=O)_2(Q_{31})$, and $-P(=O)(Q_{31})(Q_{32})$, and
- [0031]** Q_1 to Q_3 , Q_{11} to Q_{13} , Q_{21} to Q_{23} , and Q_{31} to Q_{33} may each independently be selected from hydrogen, deuterium, $-F$, $-Cl$, $-Br$, $-I$, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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 C_3 - C_{10} cycloalkyl group, a C_1 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_1 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_1 - C_{60} heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, a C_1 - C_{60} alkyl group substituted with at least one selected from deuterium, $-F$, and a cyano group, a C_6 - C_{60} aryl group substituted with at least one selected from deuterium, $-F$, and a cyano group, a biphenyl group, and a terphenyl group.
- [0032]** According to one or more embodiments, an organic light-emitting device includes: a first electrode; a second electrode facing the first electrode; and an organic layer between the first electrode and the second electrode, the organic layer including an emission layer, wherein the organic layer includes at least one of the organometallic compound described above.

BRIEF DESCRIPTION OF THE DRAWING

[0033] These and/or other aspects will become apparent and more readily appreciated from the following description of the embodiments, taken in conjunction with the FIGURE which is a schematic view of an organic light-emitting device according to an embodiment.

DETAILED DESCRIPTION

[0034] An organometallic compound according to an embodiment may be represented by Formula 1 below:



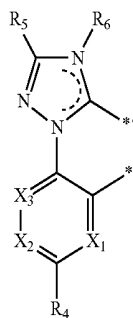
<Formula 1>

[0035] M in Formula 1 may be a third row transition metal.

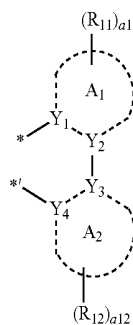
[0036] For example, M in Formula 1 may be selected from iridium (Ir), platinum (Pt), osmium (Os), titanium (Ti), zirconium (Zr), hafnium (Hf), europium (Eu), terbium (Tb), thulium (Tm), rhenium (Re), and rhodium (Rh).

[0037] In one or more embodiments, M in Formula 1 may be iridium, osmium or platinum, but embodiments of the present disclosure are not limited thereto.

[0038] In Formula 1, L_1 may be a ligand represented by Formula 2A, and n_1 may be 1, 2, or 3, wherein, when n_1 is two or more, two or more L_1 (s) may be identical to or different from each other, and L_2 may be a ligand represented by Formula 2B, and n_2 may be 0, 1, or 2, wherein, when n_2 is two or more, two or more L_2 (s) may be identical to or different from each other.



<Formula 2A>



<Formula 2B>

[0039] The sum of n_1 and n_2 in Formula 1 may be two or three.

[0040] The organometallic compound represented by Formula 1 does not have a salt form consisting of an anion and a cation and is neutral.

[0041] In one or more embodiments, M in Formula 1 may be iridium or osmium and the sum of n_1 and n_2 may be three; or M in Formula 1 may be platinum and the sum of n_1 and n_2 may be two, but embodiments of the present disclosure are not limited thereto.

[0042] * and *' in Formulae 2A and 2B each indicate a binding site to M in Formula 1.

[0043] In Formula 2A, X_1 may be N or C(R_1), X_2 may be N or C(R_2), and X_3 may be N or C(R_3).

[0044] In one or more embodiments, X_2 in Formula 2A may be N.

[0045] In one or more embodiments, X_1 in Formula 2A may be C(R_1), X_2 may be C(R_2), and X_3 may be C(R_3).

[0046] In one or more embodiments, in Formula 2A, X_1 may be C(R_1), X_2 may be N, and X_3 may be C(R_3), but embodiments of the present disclosure are not limited thereto.

[0047] In one or more embodiments, in Formula 2A, X_3 may be C(R_3), and R_3 may be an electron withdrawing group. The electron withdrawing group is the same as described herein.

[0048] In Formula 2B, Y_1 may be carbon, Y_2 and Y_3 may each independently be carbon or nitrogen, Y_4 may be nitrogen, a bond between Y_1 and Y_2 may be a single bond or a double bond, a bond between Y_2 and Y_3 may be a single bond, and a bond between Y_3 and Y_4 may be a single bond or a double bond.

[0049] For example, in Formula 2B, Y_1 to Y_3 may each be carbon, and Y_4 may be nitrogen.

[0050] Ring A_1 and ring A_2 in Formula 2B may each be a C_2 - C_{60} heterocyclic group including at least one N as a ring-forming atom.

[0051] For example, ring A_1 and ring A_2 in Formula 2B may each independently be a pyridine group, a pyrimidine group, a pyrazine group, a pyridazine group, a triazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, a phenanthroline group, a pyrazole group, an imidazole group, a triazole group, an oxazole group, an iso-oxazole group, a thiazole group, an isothiazole group, an oxadiazole group, a thiadiazol group, a benzopyrazole group, a benzimidazole group, a benzoxazole group, a benzothiazole group, a benzoxadiazole group, a benzothiadiazol group, a 5,6,7,8-tetrahydroisoquinoline group, or a 5,6,7,8-tetrahydroquinoline group.

[0052] In one or more embodiments, ring A_1 and ring A_2 in Formula 2B may each independently be a pyridine group, a pyrimidine group, a pyrazine group, a pyridazine group, a triazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, a phenanthroline group, a 5,6,7,8-tetrahydroisoquinoline group, or a 5,6,7,8-tetrahydroquinoline group.

[0053] In one or more embodiments, in Formula 2B, ring A_1 may be a pyridine group, ring A_2 may be a pyridine group, a pyrimidine group, a triazine group, a quinoline group, or an isoquinoline group, but embodiments of the present disclosure are not limited thereto.

[0054] R_1 to R_3 , R_5 , R_6 , R_{11} , and R_{12} in Formulae 2A and 2B may each independently be selected from hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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)

(Q₃), —N(Q₁)(Q₂), —B(Q₁)(Q₂), —C(=O)(Q₁), —S(=O)₂(Q₁), and —P(=O)(Q₁)(Q₂). Q₁ to Q₃ are the same as described above.

[0055] For example, R₁ to R₃, R₅, R₆, R₁₁, and R₁₂ in Formulae 2A and 2B may each independently be selected from:

[0056] hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, and a C₁-C₂₀ alkoxy group;

[0057] a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, and a C₁-C₂₀ alkoxy group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, —CD₃, —CD₂H, —CDH₂, —CF₃, —CF₂H, —CFH₂, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a pyridinyl group, a pyrimidinyl group, —Si(Q₃₁)(Q₃₂)(Q₃₃), —N(Q₃₁)(Q₃₂), —B(Q₃₁)(Q₃₂), —C(=O)(Q₃₁), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂);

[0058] a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl group, a benzofuranyl group, a benzothiofenyl 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 dibenzothiofenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, —Si(Q₃₁)(Q₃₂)(Q₃₃), —N(Q₃₁)(Q₃₂), —B(Q₃₁)(Q₃₂), —C(=O)(Q₃₁), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂); and

group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl group, a benzofuranyl group, a benzothiofenyl 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 dibenzothiofenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, an imidazopyridinyl group, and an imidazopyrimidinyl group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, —CD₃, —CD₂H, —CDH₂, —CF₃, —CF₂H, —CFH₂, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, a C₁-C₂₀ alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl group, a benzofuranyl group, a benzothiofenyl 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 dibenzothiofenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, —Si(Q₃₁)(Q₃₂)(Q₃₃), —N(Q₃₁)(Q₃₂), —B(Q₃₁)(Q₃₂), —C(=O)(Q₃₁), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂); and

[0060] —Si(Q₁)(Q₂)(Q₃), —N(Q₁)(Q₂), —B(Q₁)(Q₂), —C(=O)(Q₁), —S(=O)₂(Q₁), and —P(=O)(Q₁)(Q₂), and

[0061] Q₁ to Q₃ and Q₃₁ to Q₃₃ may each independently be selected from:

[0062] hydrogen, deuterium, —F, —Cl, —Br, —I, a cyano group, a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, a C₁-C₂₀ alkoxy group, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₂₀ aryl group, a C₁-C₂₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group;

[0063] a C₁-C₂₀ alkyl group substituted with at least one selected from deuterium, —F, and a cyano group;

[0064] a C₆-C₂₀ aryl group substituted with at least one selected from deuterium, —F, and a cyano group; and

[0065] a biphenyl group and a terphenyl group.

[0059] a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl

[0066] In one or more embodiments, R_1 to R_3 , R_5 , R_6 , R_{11} , and R_{12} in Formulae 2A and 2B may each independently be selected from:

[0067] hydrogen, deuterium, —F, —Cl, —Br, —I, a cyano group, a nitro group, a C_1 - C_{20} alkyl group, a C_2 - C_{20} alkenyl group, a C_2 - C_{20} alkynyl group, and a C_1 - C_{20} alkoxy group;

[0068] a C_1 - C_{20} alkyl group, a C_2 - C_{20} alkenyl group, a C_2 - C_{20} alkynyl group, and a C_1 - C_{20} alkoxy group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, — CD_3 , — CD_2H , — CDH_2 , — CF_3 , — CF_2H , — CFH_2 , a cyano group, a nitro group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a pyridinyl group, and a pyrimidinyl group;

[0069] a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl 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 quinolynyl group, an isoquinolynyl group, a benzoquinolynyl group, a quinoxalynyl group, a quinazolynyl group, a cinnolynyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl group, a benzofuranlyl 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 dibenzofuranlyl group, a dibenzothiophenyl group, a benzocarbazolyl group, and a dibenzocarbazolyl group;

[0070] a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl 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 quinolynyl group, an isoquinolynyl group, a benzoquinolynyl group, a quinoxalynyl group, a quinazolynyl group, a cinnolynyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl group, a benzofuranlyl 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 dibenzofuranlyl group, a dibenzothiophenyl group, a benzocarbazolyl group, and a dibenzocarbazolyl group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, — CD_3 , — CD_2H , — CDH_2 , — CF_3 , — CF_2H , — CFH_2 , a cyano

group, a nitro group, a C_1 - C_{20} alkyl group, a C_2 - C_{20} alkenyl group, a C_2 - C_{20} alkynyl group, a C_1 - C_{20} alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl 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 quinolynyl group, an isoquinolynyl group, a benzoquinolynyl group, a quinoxalynyl group, a quinazolynyl group, a cinnolynyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl group, a benzofuranlyl 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 dibenzofuranlyl group, a dibenzothiophenyl group, a benzocarbazolyl group, and a dibenzocarbazolyl group; and

[0071] —Si(Q_1)(Q_2)(Q_3), —N(Q_1)(Q_2), —B(Q_1)(Q_2), —C(=O)(Q_1), —S(=O)₂(Q_1), and —P(=O)(Q_1)(Q_2), and

[0072] Q_1 to Q_3 and Q_{31} to Q_{33} are the same as described herein.

[0073] 1) R_2 and R_3 in Formula 2A may optionally be linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group, 2) R_5 and R_6 in Formula 2A may optionally be linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group, 3) two neighboring groups of a plurality of R_{11} (s) in Formula 2B may optionally be linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group, 4) two neighboring groups of a plurality of R_{12} (s) in Formula 2B may optionally be linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group, and 5) neighboring R_{11} and R_{12} in Formula 2B may optionally be linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group. Substituents of the substituted C_4 - C_{60} carbocyclic group and the substituted C_2 - C_{60} heterocyclic group are the same as described in connection with R_{11} , the C_4 - C_{60} carbocyclic group may be, for example, a cyclopentane group, a cyclohexane group, an adamantane group, a norbornane group, a benzene group, or a naphthalene group, and the C_2 - C_{60} heterocyclic group may be, for example, a thiophene group, a furan group, a pyrrole group, a benzothiophene group, a benzofuran group, an indole group, or a pyridine group, but embodiments of the present disclosure are not limited thereto.

[0074] R_4 in Formula 2A may be an electron withdrawing group. The electron withdrawing group is the same as described herein.

[0075] a11 and a12 in Formula 2B may each independently be an integer from 0 to 10. a11 indicates the number of R_{11} (s), wherein, when a11 is two or more, two or more R_{11} (s) may be identical to or different from each other. a12

indicates the number of $R_{12}(s)$, wherein, when a_{12} is two or more, two or more $R_{12}(s)$ may be identical to or different from each other.

[0076] In one or more embodiments, in Formula 2B, the sum of a_{11} and a_{12} may be equal to or greater than one, and at least one selected from R_{11} and R_{12} may be an electron withdrawing group. The electron withdrawing group is the same as described herein.

[0077] In one or more embodiments, a_{11} in Formula 2B may be two or more, and two or more $R_{11}(s)$ may each be an electron withdrawing group.

[0078] The electron withdrawing group may be $-F$, $-Cl$, $-Br$, $-I$, a cyano group, a nitro group, or $-B(Q_1)(Q_2)$, or may be a first group substituted with at least one selected from $-F$, $-Cl$, $-Br$, $-I$, $-CF_3$, $-CF_2H$, $-CFH_2$, a cyano group, a nitro group, and $-B(Q_{31})(Q_{32})$,

[0079] provided that the first group is selected from:

[0080] a C_1 - C_{20} alkyl group, a C_2 - C_{20} alkenyl group, a C_2 - C_{20} alkynyl group, and a C_1 - C_{20} alkoxy group;

[0081] a C_1 - C_{20} alkyl group, a C_2 - C_{20} alkenyl group, a C_2 - C_{20} alkynyl group, and a C_1 - C_{20} alkoxy group, each substituted with at least one selected from deuterium, $-CD_3$, $-CD_2H$, $-CDH_2$, a hydroxyl group, an amidino group, a hydrazino group, a hydrazono group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a pyridinyl group, a pyrimidinyl group, $-Si(Q_{31})(Q_{32})(Q_{33})$, $-N(Q_{31})(Q_{32})$, $-C(=O)(Q_{31})$, $-S(=O)_2(Q_{31})$, and $-P(=O)(Q_{31})(Q_{32})$;

[0082] a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazoliny group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl 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, an imidazopyridinyl group, and an imidazopyrimidinyl group; and

[0083] a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazoliny group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl 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, an imidazopyridinyl group, and an imidazopyrimidinyl group, each substituted with at least one selected from deuterium, $-CD_3$, $-CD_2H$, $-CDH_2$, a hydroxyl group, an amidino group, a hydrazino group, a hydrazono group, a C_1 - C_{20} alkyl group, a C_2 - C_{20} alkenyl group, a C_2 - C_{20} alkynyl group, a C_1 - C_{20} alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazoliny group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl 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, an imidazopyridinyl group, $-Si(Q_{31})(Q_{32})(Q_{33})$, $-N(Q_{31})(Q_{32})$, $-C(=O)(Q_{31})$, $-S(=O)_2(Q_{31})$, and $-P(=O)(Q_{31})(Q_{32})$, and

[0084] Q_1 to Q_3 and Q_{31} to Q_{33} may each independently be selected from:

[0085] hydrogen, deuterium, $-F$, $-Cl$, $-Br$, $-I$, a cyano group, a C_1 - C_{20} alkyl group, a C_2 - C_{20} alkenyl group, a C_2 - C_{20} alkynyl group, a C_1 - C_{20} alkoxy group, a C_3 - C_{10} cycloalkyl group, a C_1 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_1 - C_{10} heterocycloalkenyl group, a C_6 - C_{20} aryl group, a C_1 - C_{20} heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group;

[0086] a C_1 - C_{20} alkyl group substituted with at least one selected from deuterium, $-F$, and a cyano group;

[0087] a C_6 - C_{20} aryl group substituted with at least one selected from deuterium, $-F$, and a cyano group; and

[0088] a biphenyl group and a terphenyl group.

[0089] For example, it is understood from the present disclosure that, 1) when the electron withdrawing group is a first group substituted with at least one —F and the first group is a “methyl group”, the electron withdrawing group is a “methyl group substituted with at least one —F (for example, —CF₃, —CHF₂, or —CH₂F)”, 2) when the electron withdrawing group is a first group substituted with at least one —F and the first group is a “methyl group substituted with at least one deuterium”, the electron withdrawing group is a “methyl group substituted with at least one —F and at least one deuterium (for example, —CDF₂ or —CD₂F)”.

[0090] In one or more embodiments, the electron withdrawing group may be —F, —Cl, —Br, —I, a cyano group, a nitro group, or —B(Q₁)(Q₂), or may be a first group substituted with at least one selected from —F, —Cl, —Br, —I, —CF₃, —CF₂H, —CFH₂, a cyano group, a nitro group, and —B(Q₃₁)(Q₃₂).

[0091] provided that the first group is selected from:

[0092] a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, and a C₁-C₂₀ alkoxy group;

[0093] a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, and a C₁-C₂₀ alkoxy group, each substituted with at least one selected from deuterium, —CD₃, —CD₂H, —CDH₂, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a pyridinyl group, and a pyrimidinyl group;

[0094] a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, and a fluorenyl group; and

[0095] a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, and a fluorenyl group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, —CD₃, —CD₂H, —CDH₂, a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, a C₁-C₂₀ alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, and a fluorenyl group, and

[0096] Q₁, Q₂, Q₃₁, and Q₃₂ may each independently be selected from:

[0097] hydrogen, deuterium, —F, —Cl, —Br, —I, a cyano group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, and a naphthyl group;

[0098] a C₁-C₂₀ alkyl group substituted with at least one selected from deuterium, —F, and a cyano group;

[0099] a phenyl group and a naphthyl group, each substituted with at least one selected from deuterium, —F, and a cyano group; and

[0100] a biphenyl group and a terphenyl group.

[0101] In one or more embodiments, the electron withdrawing group may be —F, —Cl, —Br, —I, a cyano group, a nitro group, or —B(Q₁)(Q₂), or may be a first group selected from at least one selected from —F, —Cl, —Br, —I, —CF₃, —CF₂H, —CFH₂, a cyano group, a nitro group, and —B(Q₃₁)(Q₃₂),

[0102] provided that the first group is selected from:

[0103] a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, and a C₁-C₂₀ alkoxy group; and

[0104] a C₁-C₂₀ alkyl group and a C₁-C₂₀ alkoxy group, each substituted with at least one selected from deuterium, —CD₃, —CD₂H, —CDH₂, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a pyridinyl group, and a pyrimidinyl group, and

[0105] Q₁, Q₂, Q₃₁ and Q₃₂ may each independently be selected from:

[0106] hydrogen, deuterium, —F, —Cl, —Br, —I, a cyano group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, and a naphthyl group;

[0107] a C₁-C₂₀ alkyl group substituted with at least one selected from deuterium, —F, and a cyano group;

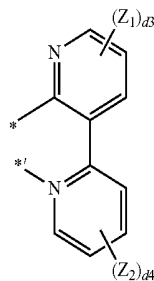
[0108] a phenyl group and a naphthyl group, each substituted with at least one selected from deuterium, —F, and a cyano group; and

[0109] a biphenyl group and a terphenyl group,

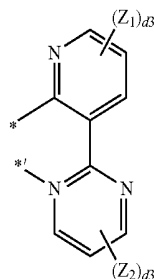
[0110] but embodiments of the present disclosure are not limited thereto.

[0111] L₂ may be a ligand selected from groups represented by Formulae 2B-1 to 2B-60:

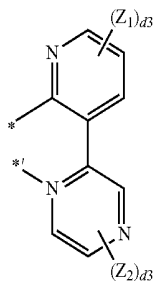
Formula 2B-1



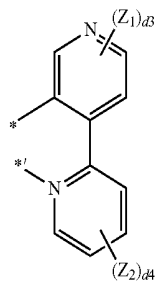
Formula 2B-2



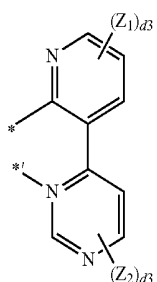
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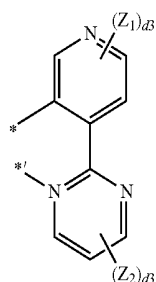
Formula 2B-3



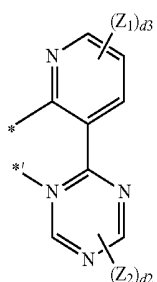
Formula 2B-8



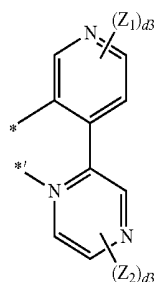
Formula 2B-4



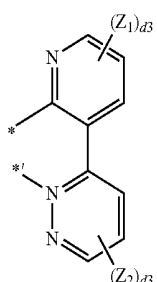
Formula 2B-9



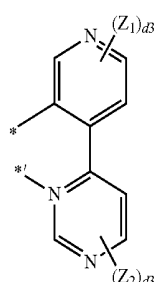
Formula 2B-5



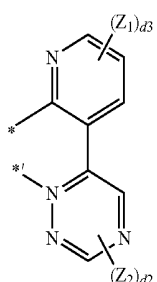
Formula 2B-10



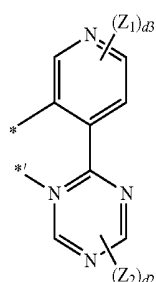
Formula 2B-6



Formula 2B-11

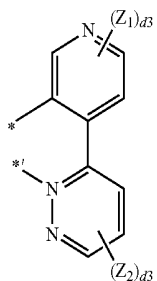


Formula 2B-7

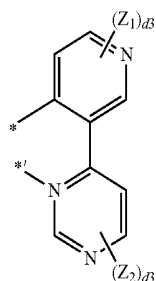


Formula 2B-12

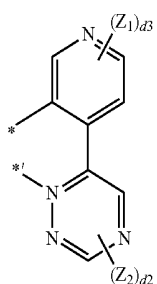
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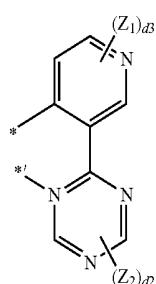
Formula 2B-13



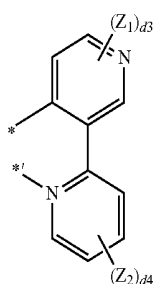
Formula 2B-18



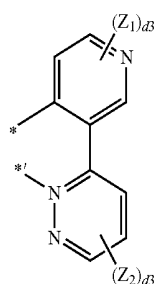
Formula 2B-14



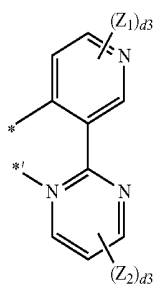
Formula 2B-19



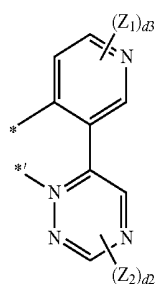
Formula 2B-15



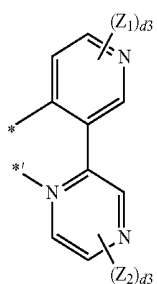
Formula 2B-20



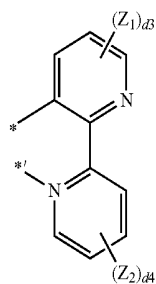
Formula 2B-16



Formula 2B-21

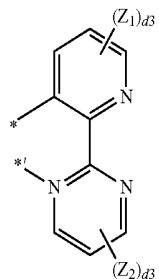


Formula 2B-17



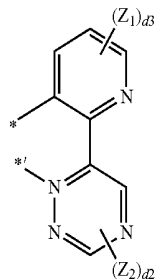
Formula 2B-22

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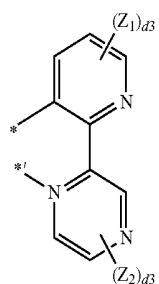


Formula 2B-23

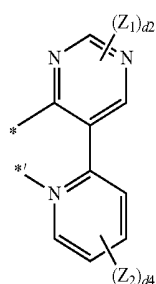
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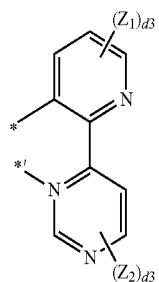
Formula 2B-28



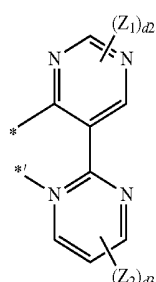
Formula 2B-24



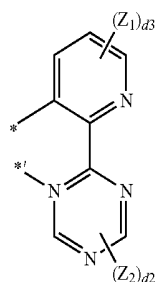
Formula 2B-29



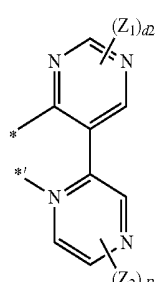
Formula 2B-25



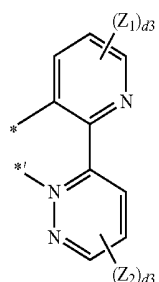
Formula 2B-30



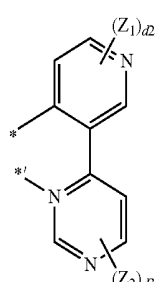
Formula 2B-26



Formula 2B-31

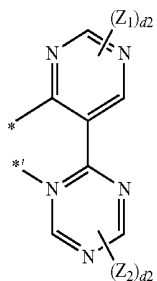


Formula 2B-27

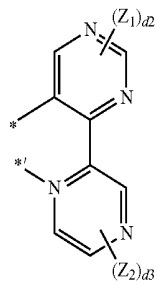


Formula 2B-32

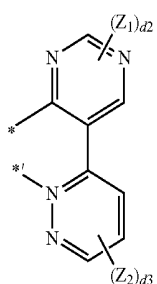
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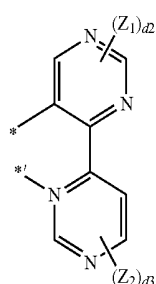
Formula 2B-33



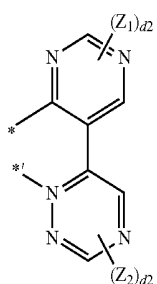
Formula 2B-38



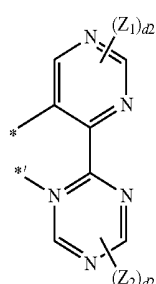
Formula 2B-34



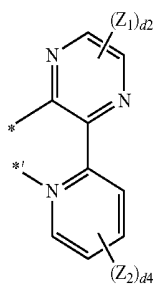
Formula 2B-39



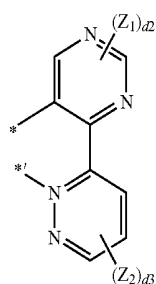
Formula 2B-35



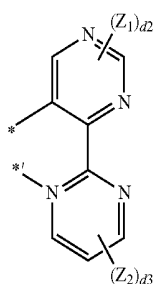
Formula 2B-40



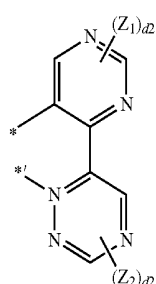
Formula 2B-36



Formula 2B-41

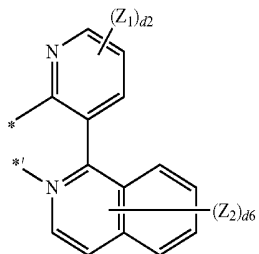


Formula 2B-37

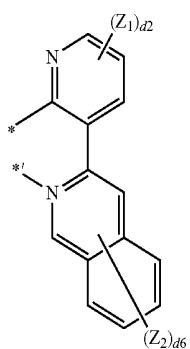


Formula 2B-42

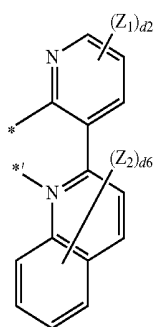
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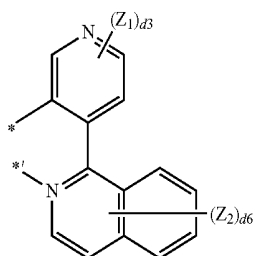
Formula 2B-43



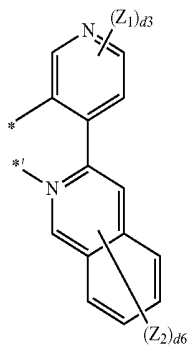
Formula 2B-44



Formula 2B-45

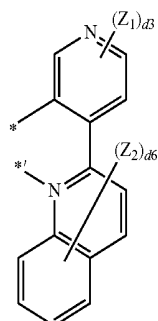


Formula 2B-46

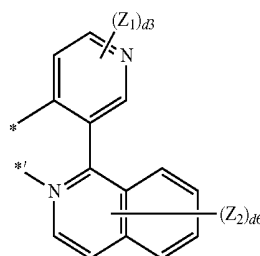


Formula 2B-47

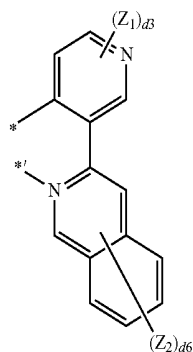
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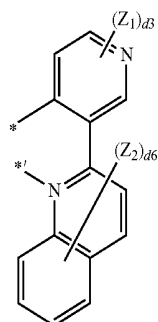
Formula 2B-48



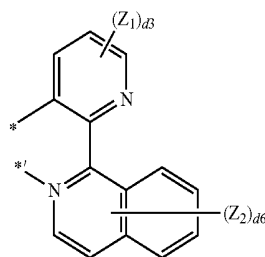
Formula 2B-49



Formula 2B-50

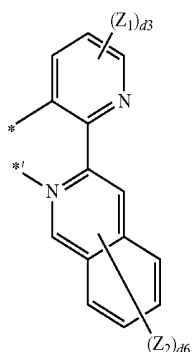


Formula 2B-51

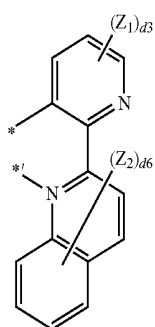


Formula 2B-52

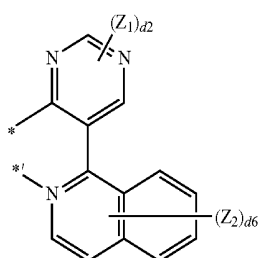
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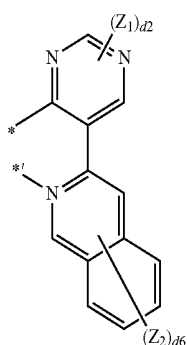
Formula 2B-53



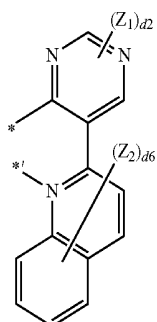
Formula 2B-54



Formula 2B-55

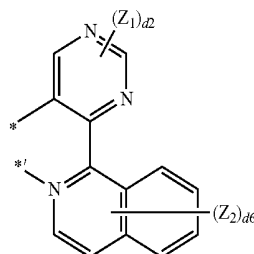


Formula 2B-56

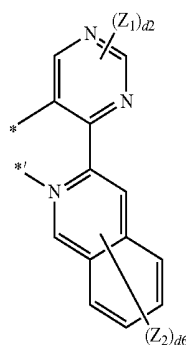


Formula 2B-57

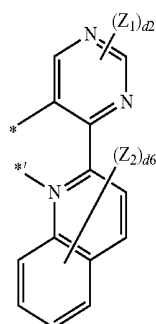
-continued



Formula 2B-58



Formula 2B-59



Formula 2B-60

[0112] In Formulae 2B-1 to 2-60,

[0113] Z_1 and Z_2 are the same as described in connection with R_{11} and R_{12} , respectively,

[0114] d_2 may be an integer from 0 to 2,

[0115] d_3 may be an integer from 0 to 3,

[0116] d_4 may be an integer from 0 to 4,

[0117] d_6 may be an integer from 0 to 6, and

[0118] * and *' each indicate a binding site to M in Formula 1.

[0119] In one or more embodiments, Formulae 2B-1 to 2B-60 have at least one Z_1 and at least one Z_2 , provided that at least one selected from Z_1 and Z_2 is an electron withdrawing group. The electron withdrawing group is the same as described herein.

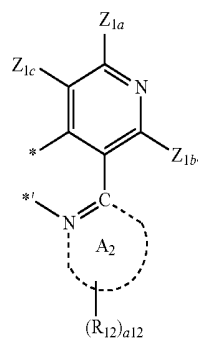
[0120] In one or more embodiments, Formulae 2B-1 to 2B-60 have at least one Z_1 , provided that at least one of Z_1 (s) is an electron withdrawing group. The electron withdrawing group is the same as described herein.

[0121] For example, L_2 in Formula 1 may be a ligand selected from groups represented by Formula 2B-A:

-continued

<Formula 2B-A>

[0122]



[0123] In Formula 2B-A,

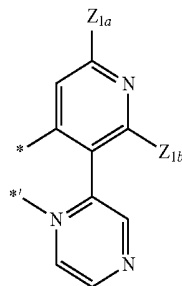
[0124] Z_{1a} , Z_{1b} , and Z_{1c} are each independently the same as described in connection with R_{11} ,

[0125] ring A_2 , R_{12} , and a_{12} are each independently the same as described above, and

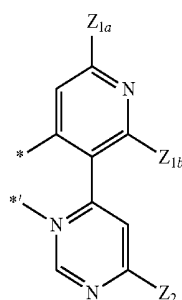
[0126] * and *' each indicate a binding site to M in Formula 1.

[0127] In one or more embodiments, L_2 in Formula 1 may be a ligand selected from groups represented by Formulae 2B(1) to 2B(8):

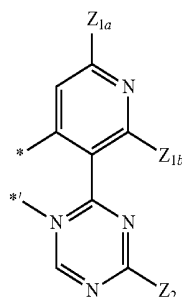
Formula 2B(3)



Formula 2B(4)

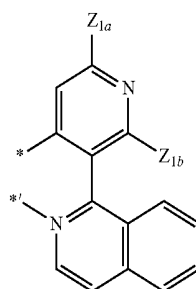
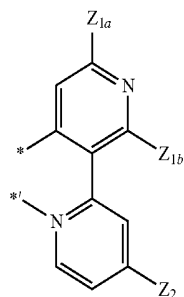


Formula 2B(5)



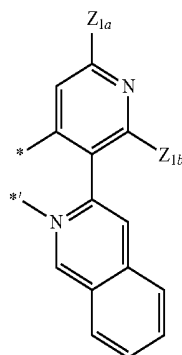
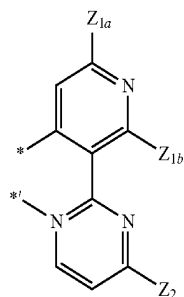
Formula 2B(6)

Formula 2B(1)



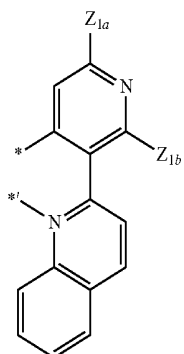
Formula 2B(7)

Formula 2B(2)



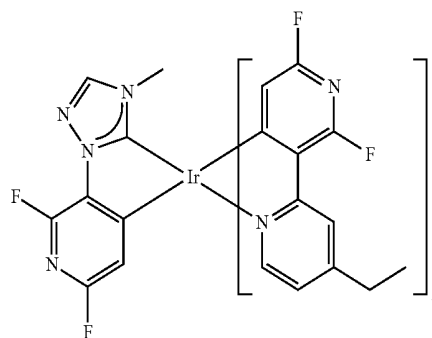
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Formula 2B(8)



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4

[0128] In Formulae 2B(1) and 2B(8),

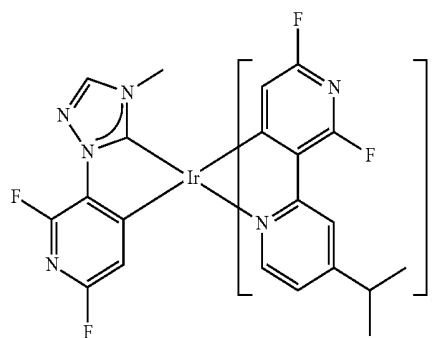
[0129] Z_{1a} and Z_{1b} are the same as described in connection with R_{11} ,

[0130] Z_2 is the same as described in connection with R_{12} , and

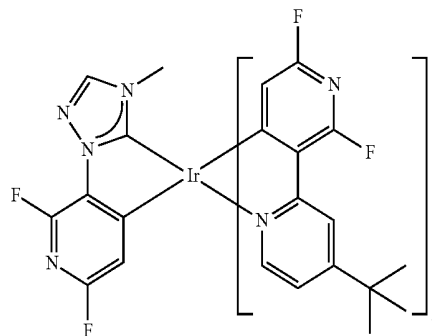
[0131] * and *' each indicate a binding site to M in Formula 1.

[0132] In one or more embodiments, at least one selected from Z_{1a} and Z_{1b} in Formulae 2B-A and 2B(1) to 2B(8) may be an electron withdrawing group, or both Z_{1a} and Z_{1b} may be an electron withdrawing group at the same time, but embodiments of the present disclosure are not limited thereto. The electron withdrawing group is the same as described herein.

[0133] In one or more embodiments, the organometallic compound may be selected from Compounds 1 to 96, but embodiments of the present disclosure are not limited thereto:

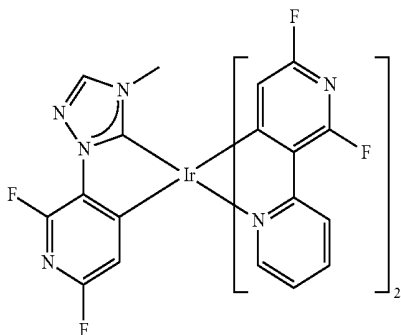


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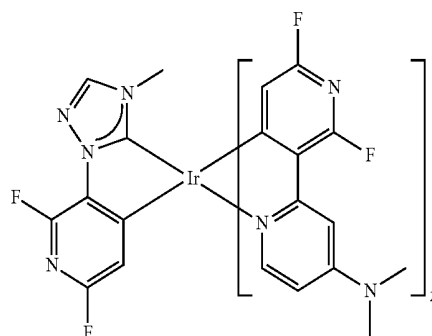


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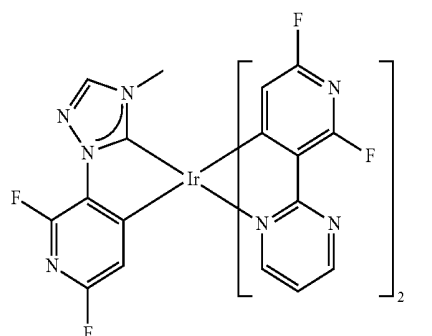
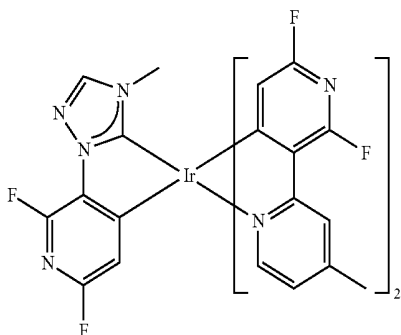
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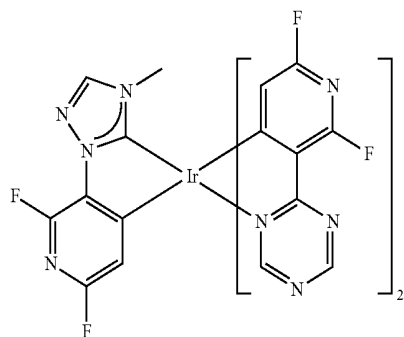
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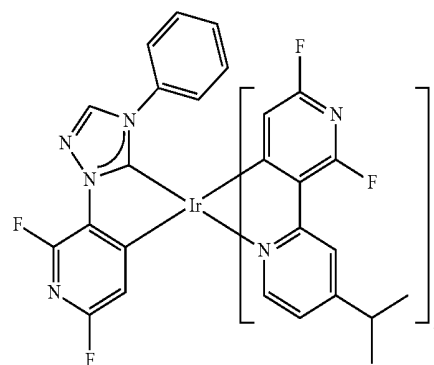


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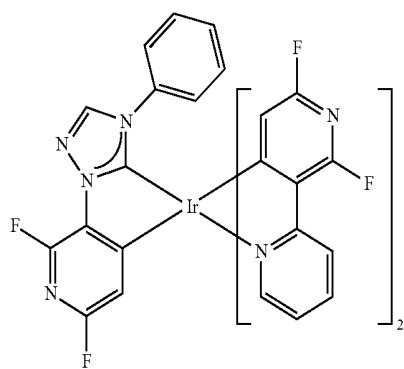


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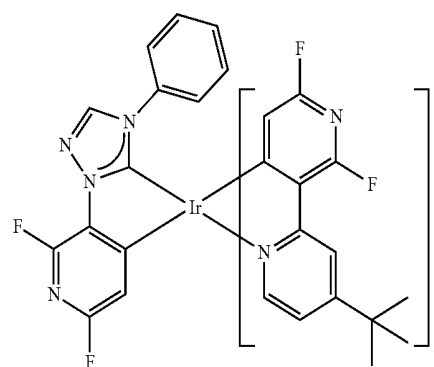
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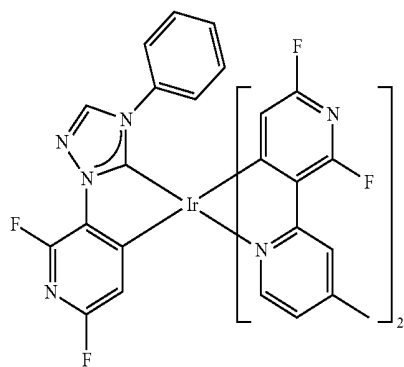
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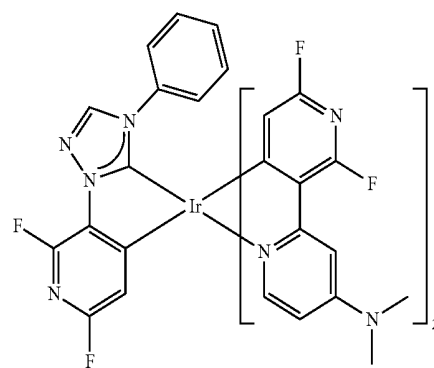
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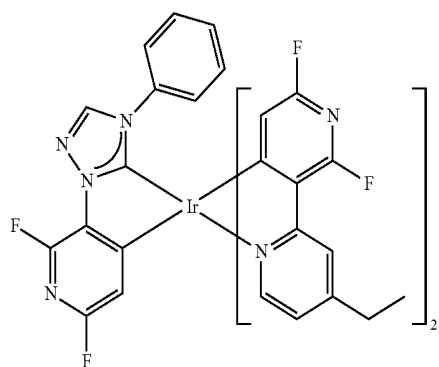
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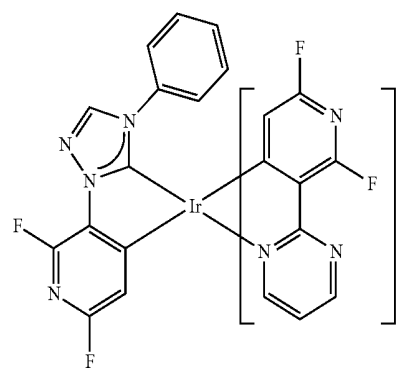
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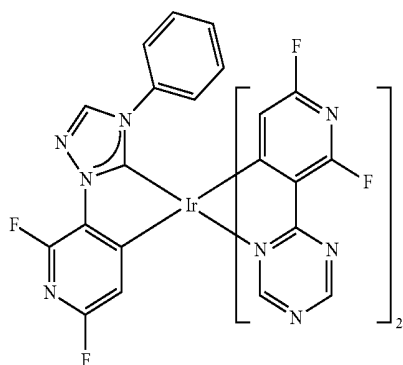


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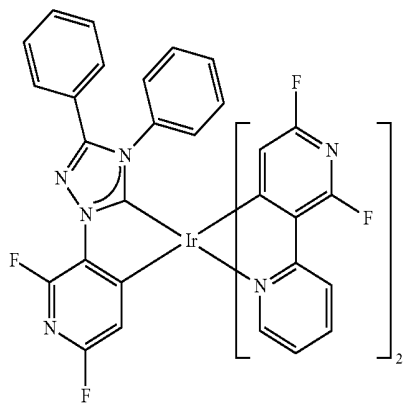


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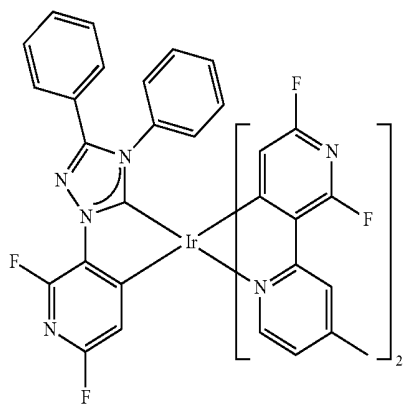
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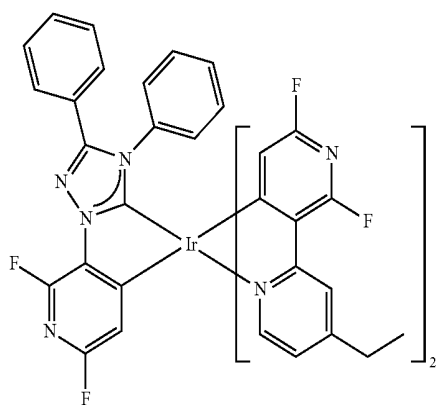
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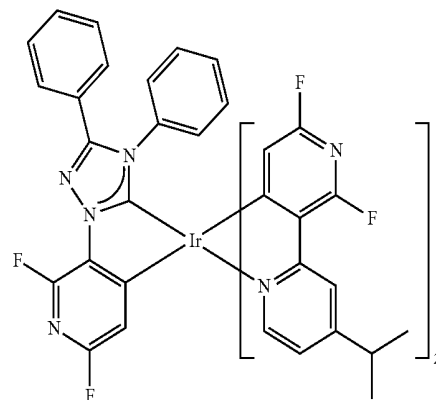


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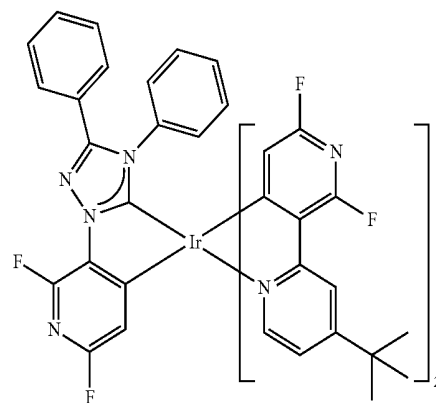


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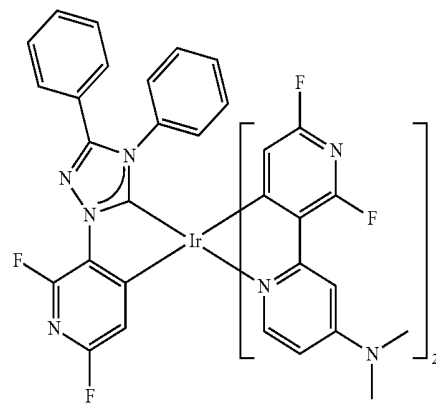
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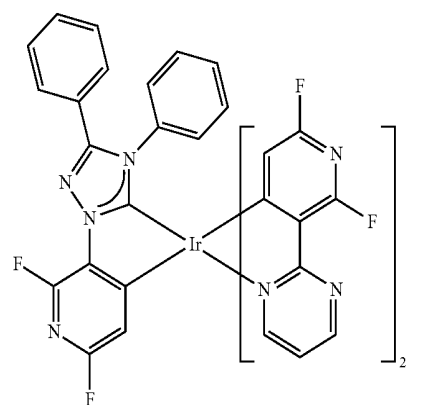
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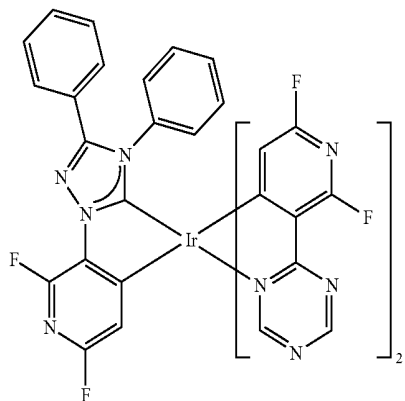


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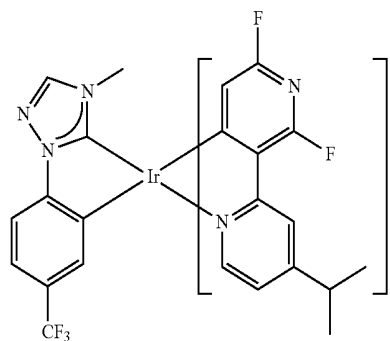
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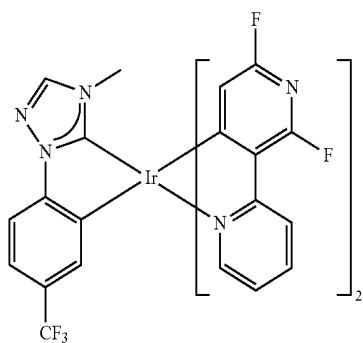


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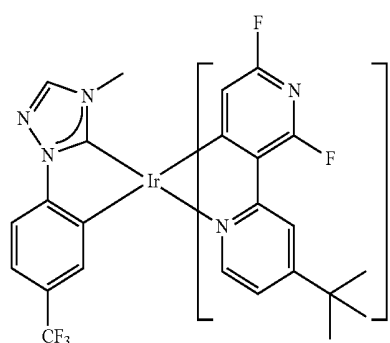
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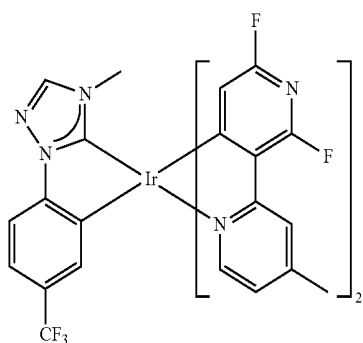
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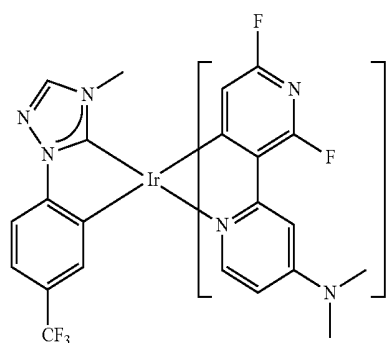
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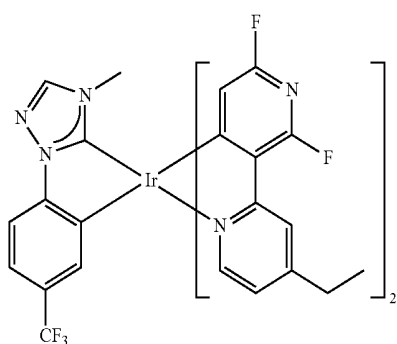
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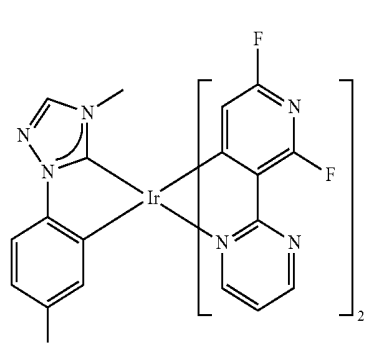
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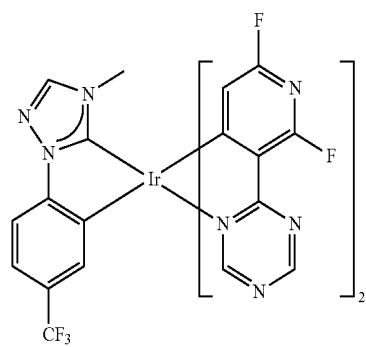


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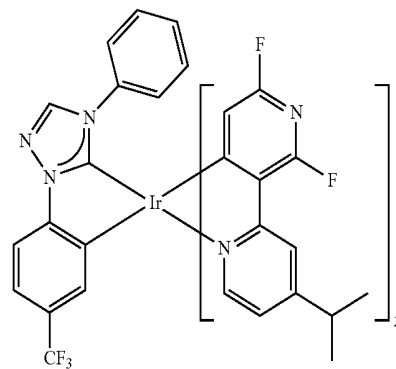
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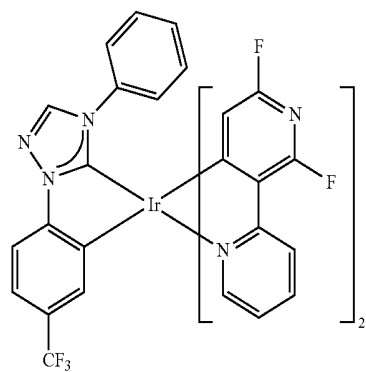
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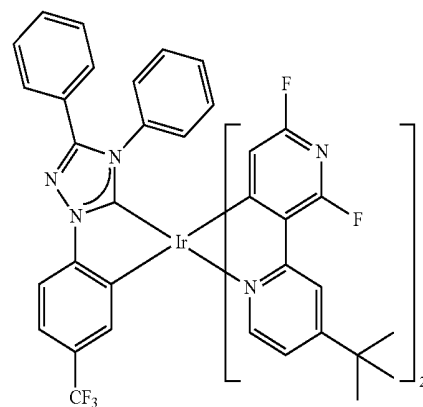


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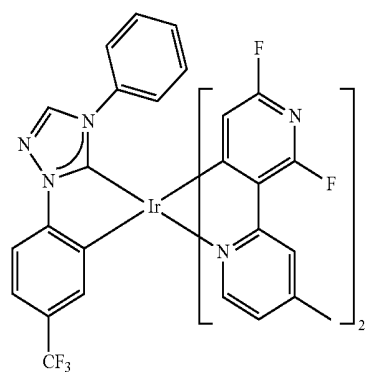


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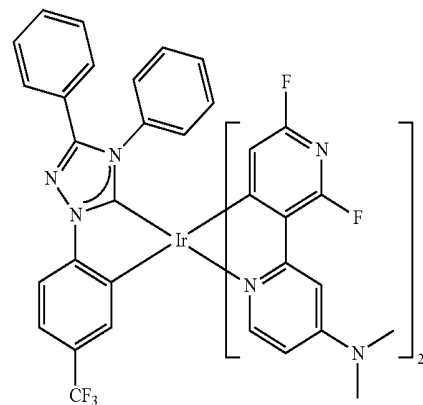


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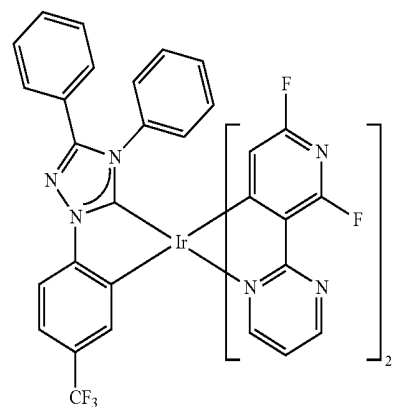
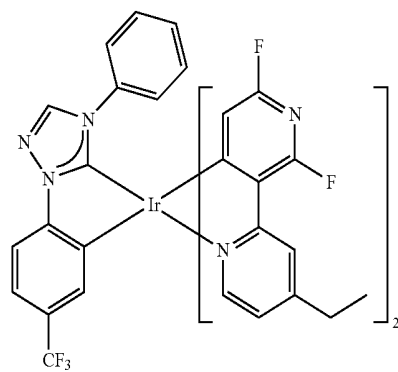
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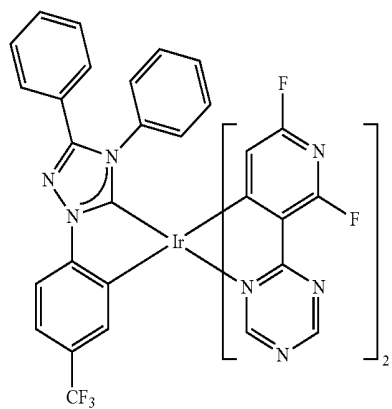
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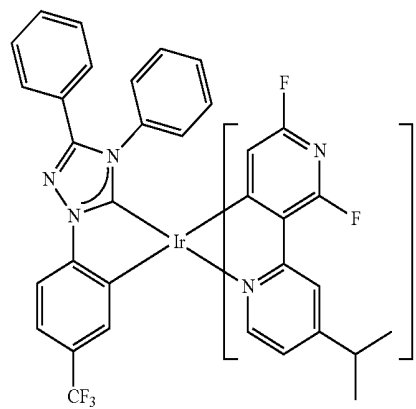


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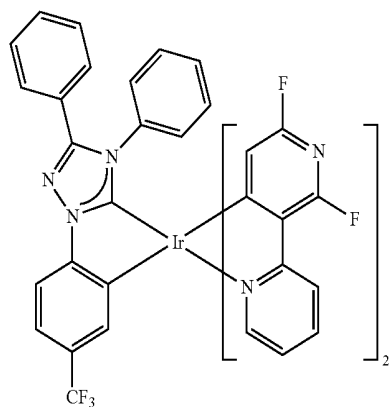


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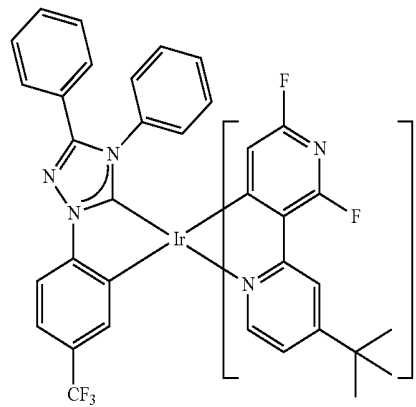
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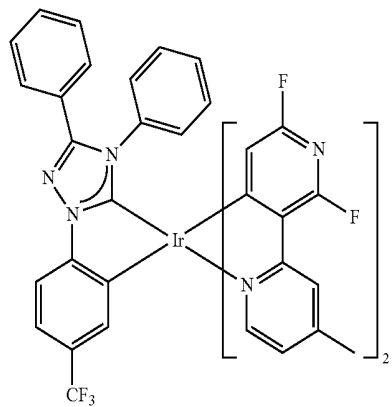
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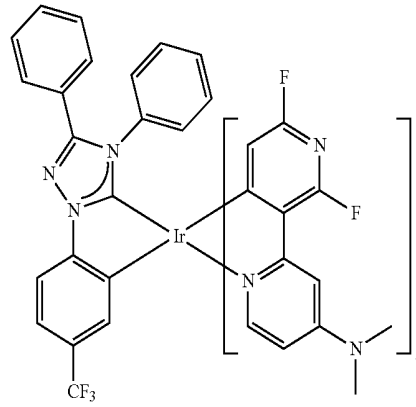
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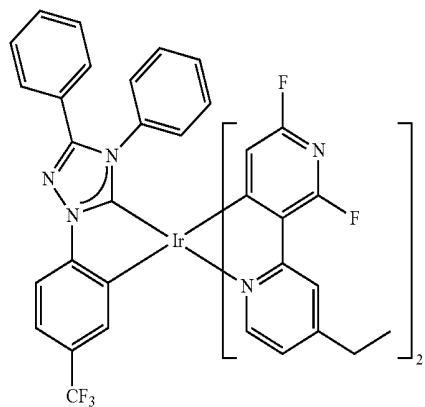
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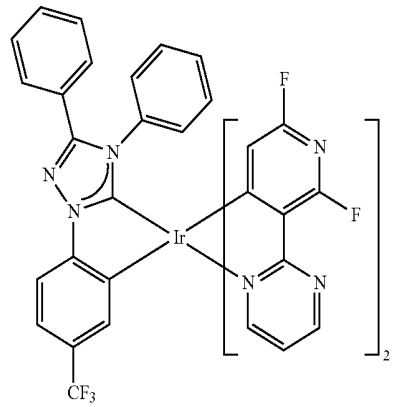
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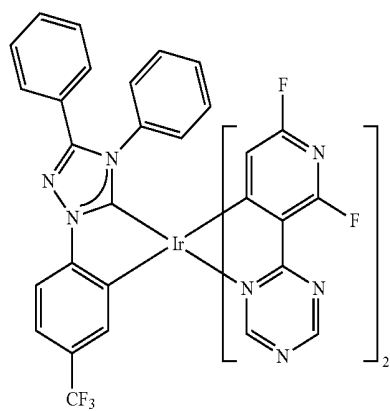


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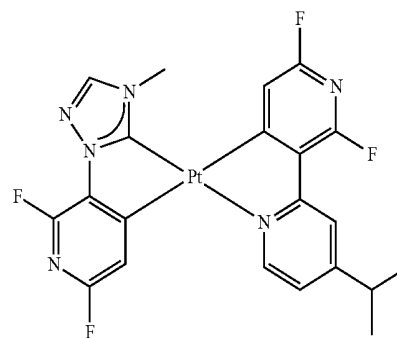
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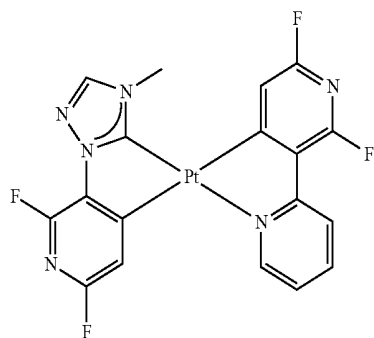
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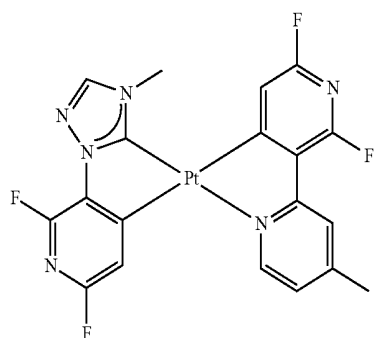
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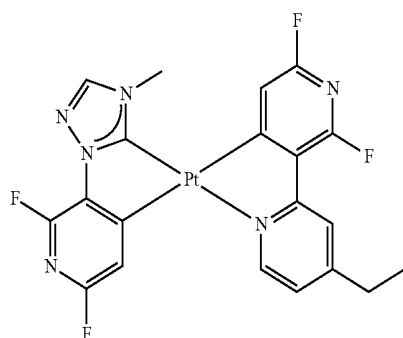
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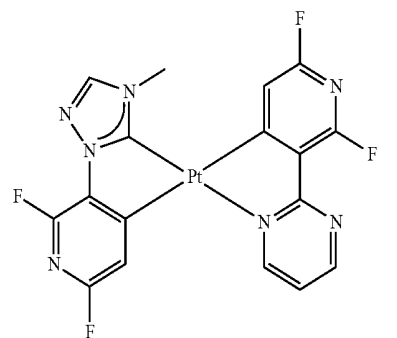
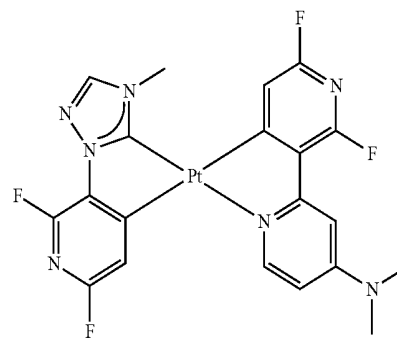


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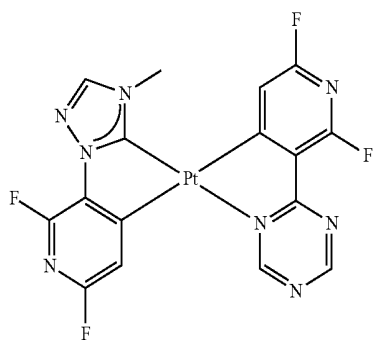
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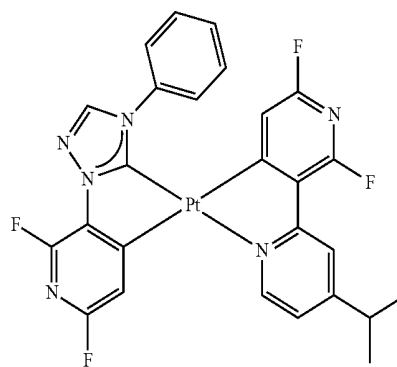


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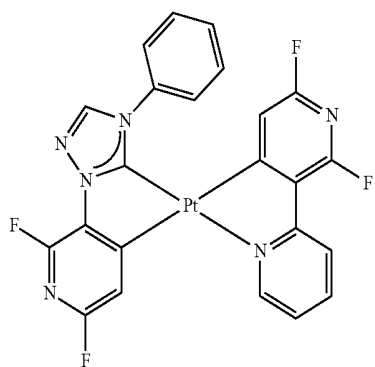


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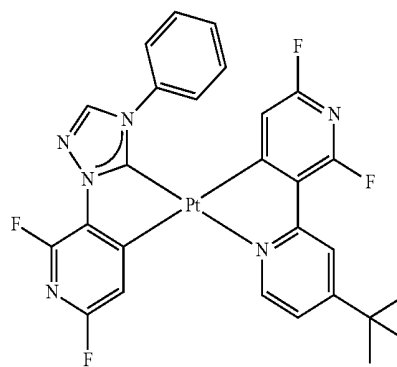
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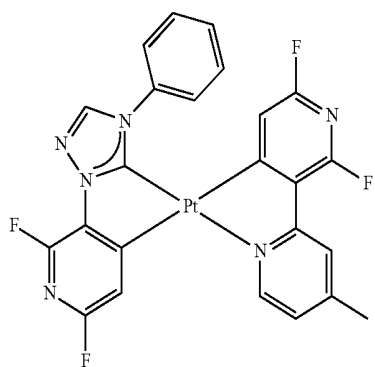
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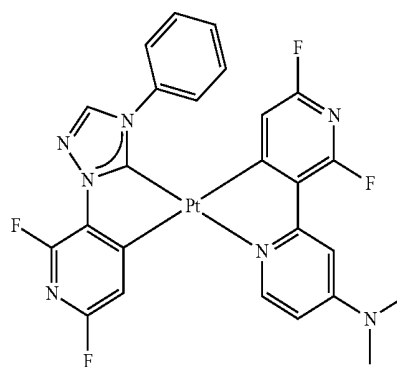
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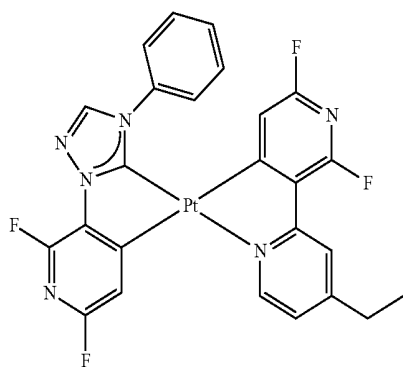
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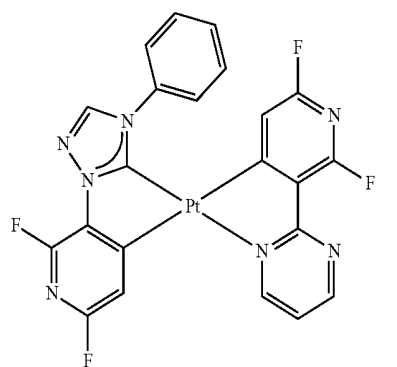
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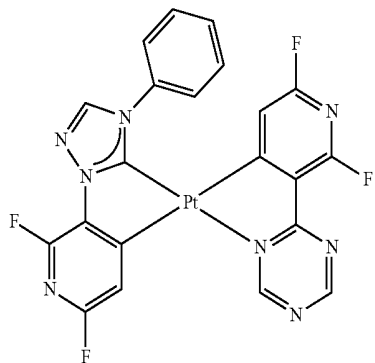


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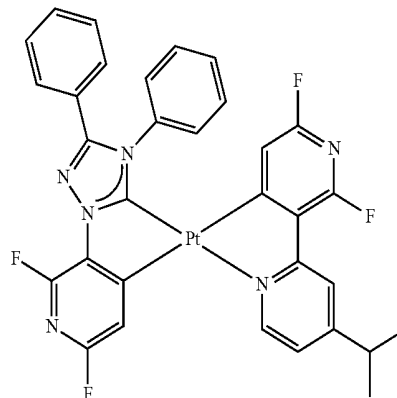
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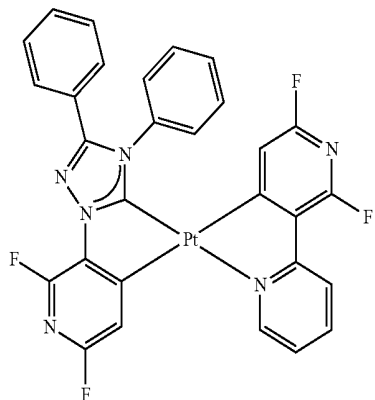
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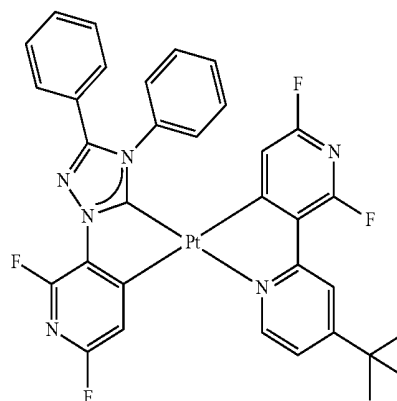


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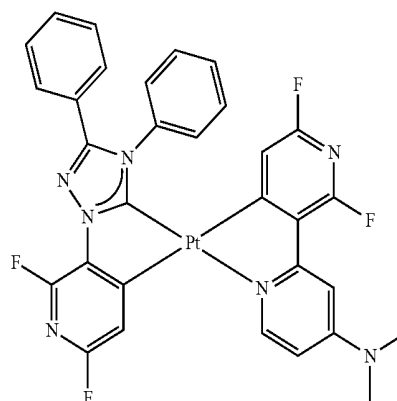
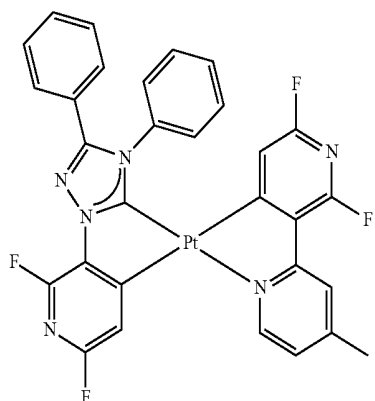


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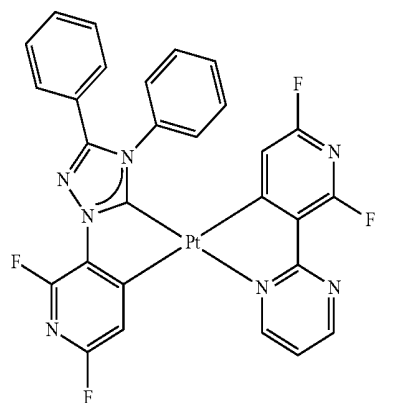
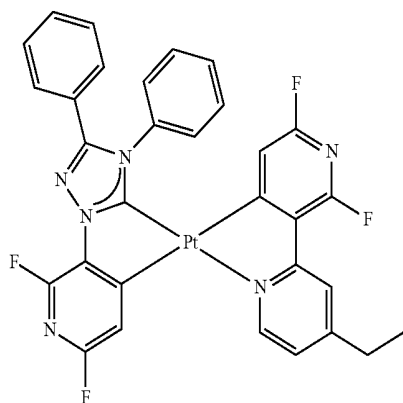
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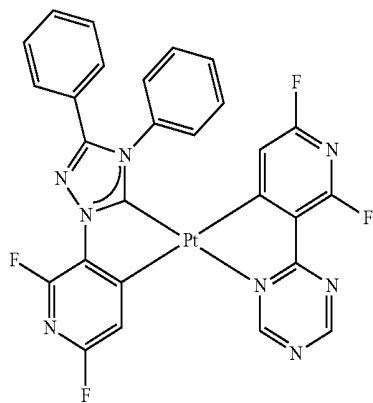


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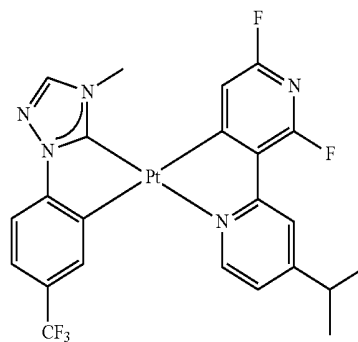


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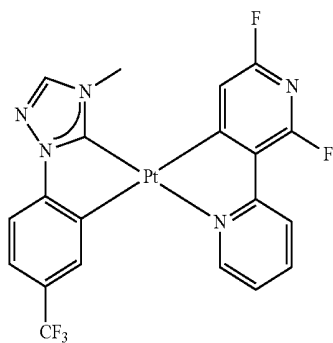
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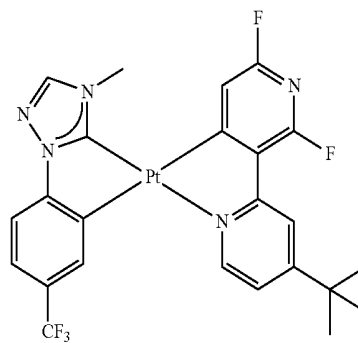


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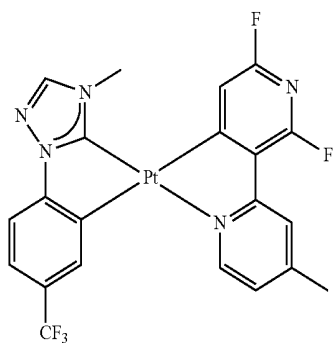
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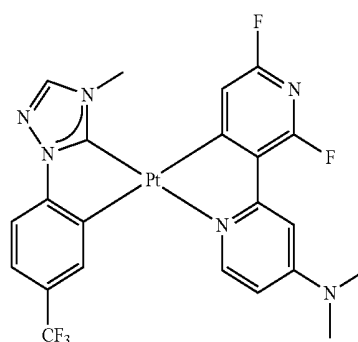
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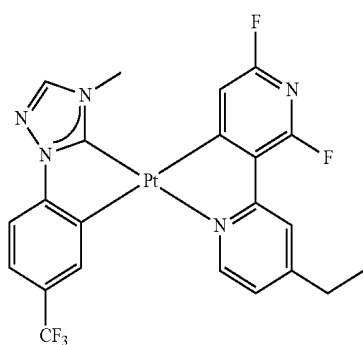
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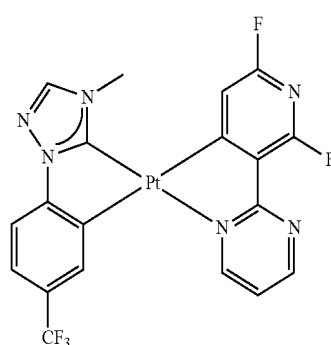
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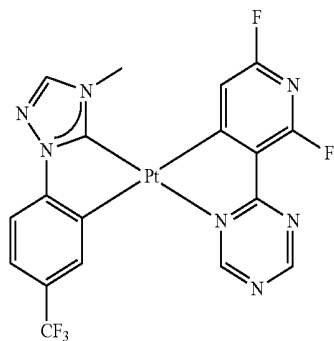
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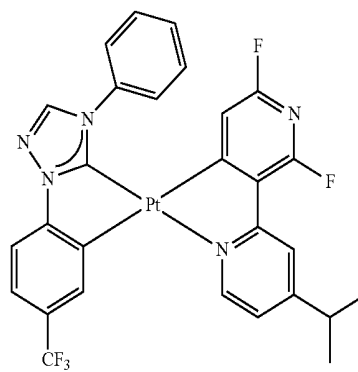


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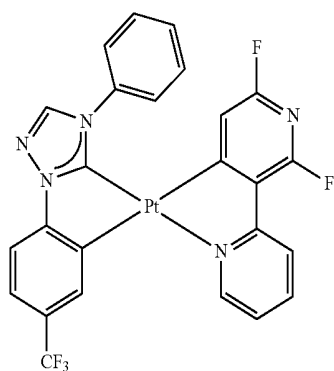


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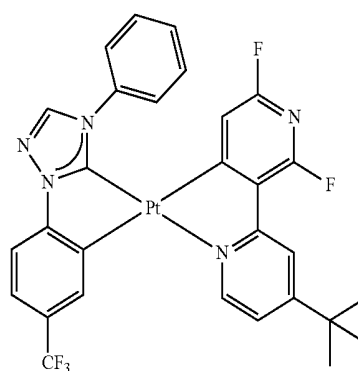
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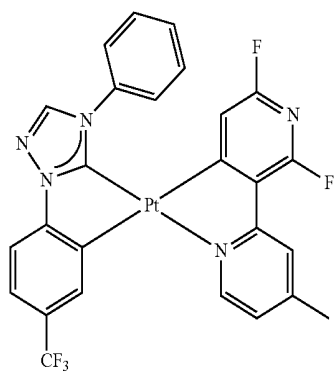
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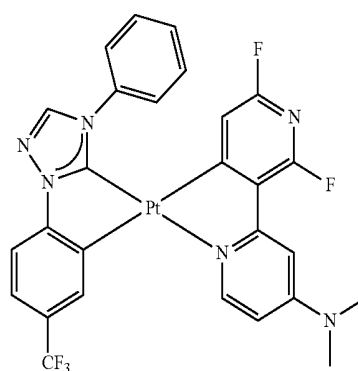
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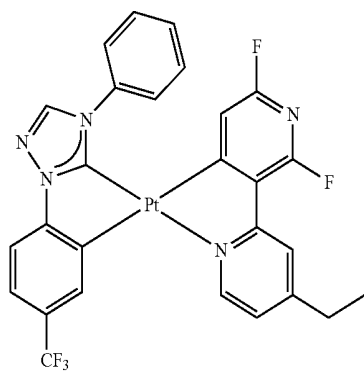
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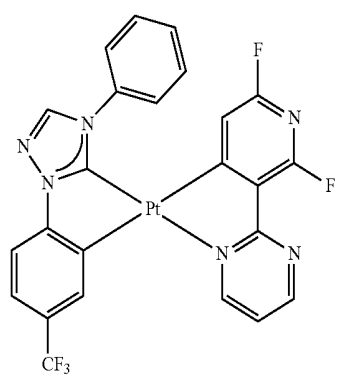
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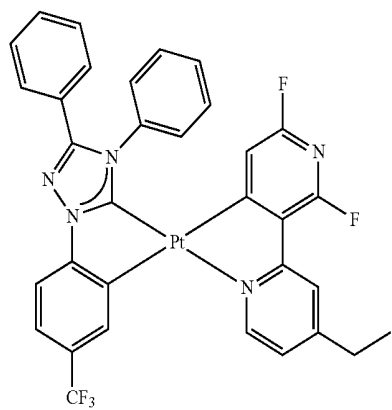
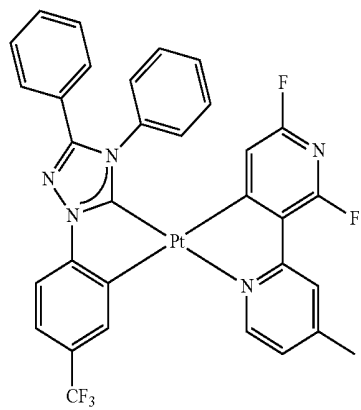
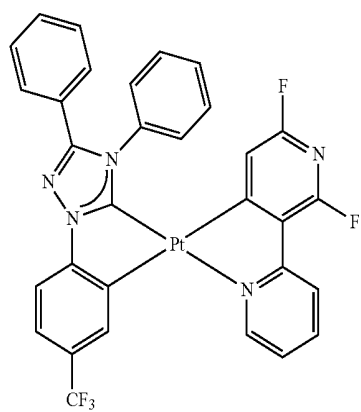
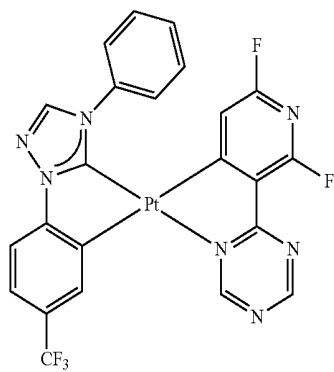


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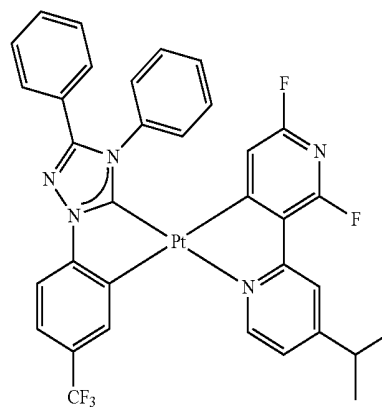
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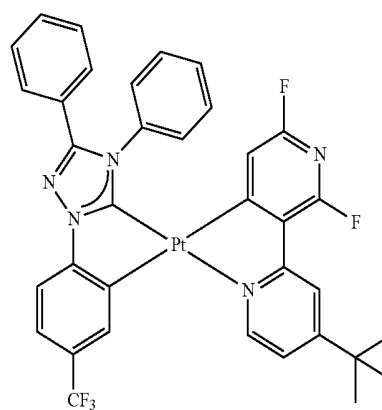
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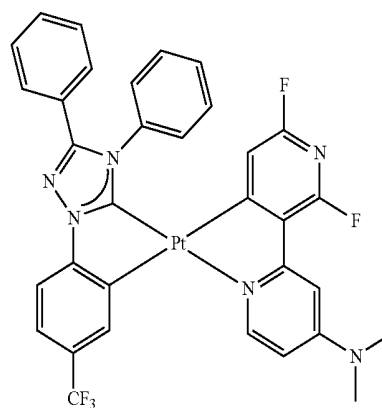
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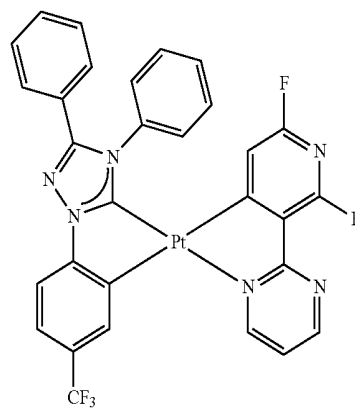
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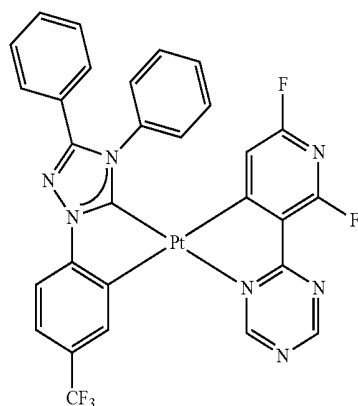
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[0134] The organometallic compound represented by Formula 1 essentially includes at least one ligand represented by Formula 2A. Since Formula 2A is a carbene ligand including a 1,2,4-triazole ring and R_4 is essentially an electron withdrawing group, the organometallic compound has a relatively high lowest unoccupied molecular orbital (LUMO) energy level and a relatively low highest occupied molecular orbital (HOMO) energy level. Accordingly, the organometallic compound may have a large energy bandgap, a ligand to ligand charge transfer transition (LLCT) is minimized to thereby emit deep blue light.

[0135] Also, the organometallic compound represented by Formula 1 may optionally further include a ligand represented by Formula 2B. In Formula 2B, Y_1 may be carbon, Y_4 may be nitrogen, and ring A_1 and ring A_2 may each be a C_2-C_{60} heterocyclic group including at least one N as a ring-forming atom. Thus, a LUMO energy level of the organometallic compound may be relatively further increased. Accordingly, the organometallic compound may emit deep blue light having high color purity.

[0136] For example, the organometallic compound may emit blue light having a maximum emission wavelength of 440 nm to 495 nm, for example, 445 nm to 465 nm, (for example, bottom-emission CIE_{x,y} color coordinates $x=0.14$, $y=0.06$ to 0.08), but embodiments of the present disclosure are not limited thereto. Accordingly, the organometallic compound represented by Formula 1 may be used to manufacture an organic light-emitting device that emits deep blue light.

[0137] Synthesis methods of the organometallic compound represented by Formula 1 may be recognizable by those of ordinary skill in the art by referring to Examples provided below.

[0138] At least one of the organometallic compound represented by Formula 1 may be used between a pair of electrodes of an organic light-emitting device. For example, the organometallic compound may be included in an emission layer. The organometallic compound included in the emission layer may act as a dopant.

[0139] Accordingly, provided is an organic light-emitting device including: a first electrode; a second electrode facing the first electrode; and an organic layer between the first electrode and the second electrode, the organic layer including an emission layer, wherein the organic layer includes at least one of the organometallic compound represented by Formula 1.

[0140] The expression “(an organic layer) includes at least one of organometallic compounds” used herein may include a case in which “(an organic layer) includes identical organometallic compounds represented by Formula 1” and a case in which “(an organic layer) includes two or more different organometallic compounds represented by Formula 1.”

[0141] For example, the organic layer may include, as the organometallic compound, only Compound 1. In this regard, Compound 1 may exist in an emission layer of the organic light-emitting device. In one or more embodiments, the organic layer may include, as the organometallic 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 all exist in an emission layer), or different layers (for example, Compound 1 may exist in an emission layer and Compound 2 may exist in an electron transport region).

[0142] In one or more embodiments,

[0143] the first electrode of the organic light-emitting device may be an anode,

[0144] the second electrode of the organic light-emitting device may be a cathode,

[0145] the organic layer may further include a hole transport region between the first electrode and the emission layer and an electron transport region between the emission layer and the second electrode,

[0146] the hole transport region may include a hole injection layer, a hole transport layer, an emission auxiliary layer, an electron blocking layer, or any combination thereof, and

[0147] the electron transport region may include a hole blocking layer, an electron transport layer, an electron injection layer, or any combination thereof.

[0148] The term “organic layer” as used herein refers to a single layer and/or a plurality of layers 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.

DESCRIPTION OF THE FIGURE

[0149] The FIG. 1 is a schematic view of an organic light-emitting device 10 according to an embodiment. The organic light-emitting device 10 includes a first electrode 110, an organic layer 150, and a second electrode 190.

[0150] Hereinafter, the structure of the organic light-emitting device 10 according to an embodiment and a method of manufacturing the organic light-emitting device 10 will be described in connection with the FIGURE.

[First Electrode 110]

[0151] In the FIGURE, 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 a plastic substrate, each having excellent mechanical strength, thermal stability, transparency, surface smoothness, ease of handling, and water resistance.

[0152] The first electrode **110** may be formed by depositing or sputtering a material for forming the first electrode **110** on the substrate. When the first electrode **110** is an anode, the material for forming a first electrode may be selected from materials with a high work function to facilitate hole injection.

[0153] The first electrode **110** may be a reflective electrode, a semi-transmissive electrode, or a transmissive electrode. When the first electrode **110** is a transmissive electrode, a material for forming a first electrode may be selected from indium tin oxide (ITO), indium zinc oxide (IZO), tin oxide (SnO₂), zinc oxide (ZnO), and any combinations thereof, but embodiments of the present disclosure are not limited thereto. When the first electrode **110** is a semi-transmissive electrode or a reflective electrode, as a material for forming the first electrode **110**, magnesium (Mg), silver (Ag), aluminum (Al), aluminum-lithium (Al—Li), calcium (Ca), magnesium-indium (Mg—In), magnesium-silver (Mg—Ag), or any combination thereof may be used. However, the material for forming the first electrode **110** is not limited thereto.

[0154] The first electrode **110** may have a single-layered structure, or a multi-layered structure including two or more layers. For example, the first electrode **110** may have a three-layered structure of ITO/Ag/ITO, but the structure of the first electrode **110** is not limited thereto.

[Organic Layer **150**]

[0155] The organic layer **150** is disposed on the first electrode **110**. The organic layer **150** may include an emission layer.

[0156] The organic layer **150** may further include a hole transport region between the first electrode **110** and the emission layer, and an electron transport region between the emission layer and the second electrode **190**.

[Hole Transport Region in Organic Layer **150**]

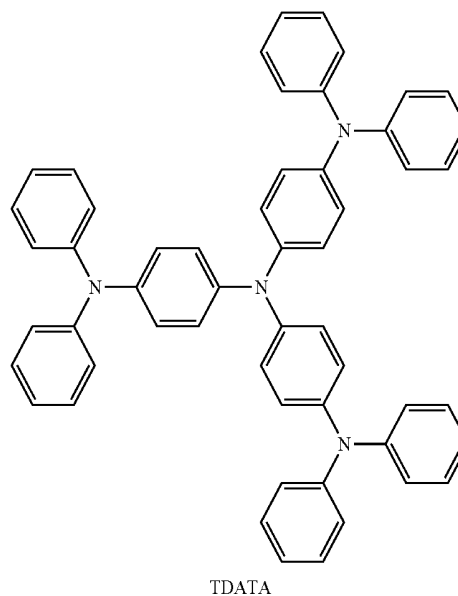
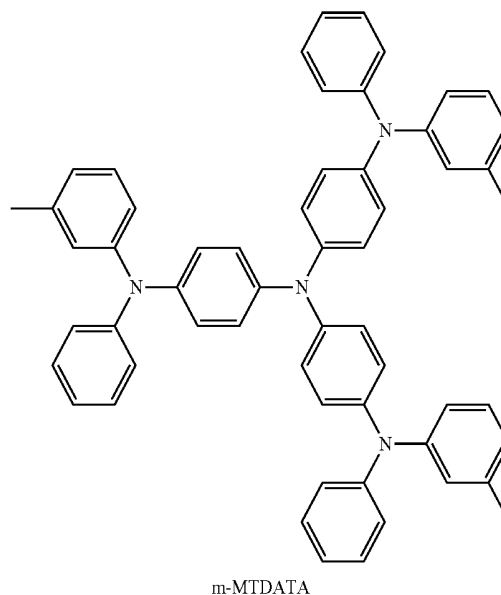
[0157] The hole transport region may have i) a single-layered structure including a single layer including a single material, ii) a single-layered structure including a single layer including a plurality of different materials, or iii) a multi-layered structure having a plurality of layers including a plurality of different materials.

[0158] The hole transport region may include at least one layer selected from a hole injection layer (HIL), a hole transport layer (HTL), an emission auxiliary layer, and an electron blocking layer (EBL).

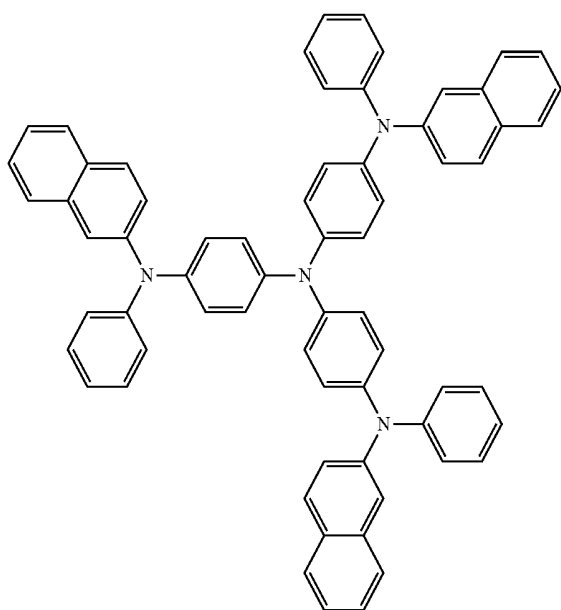
[0159] For example, the hole transport region may have a single-layered structure including a single layer including a plurality of different materials, or a multi-layered structure having a hole injection layer/hole transport layer structure, a hole injection layer/hole transport layer/emission auxiliary layer structure, a hole injection layer/emission auxiliary layer structure, a hole transport layer/emission auxiliary layer structure, or a hole injection layer/hole transport layer/electron blocking layer structure, wherein for each structure, constituting layers are sequentially stacked from the first

electrode **110** in this stated order, but the structure of the hole transport region is not limited thereto.

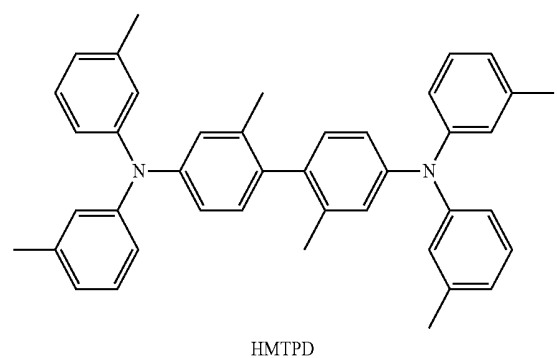
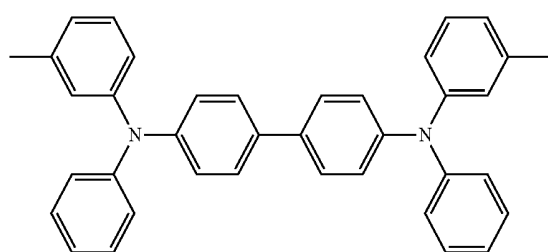
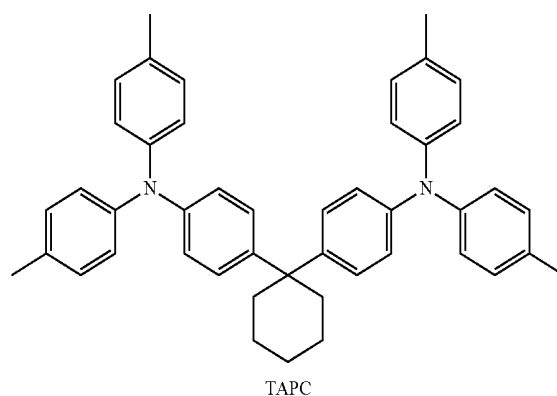
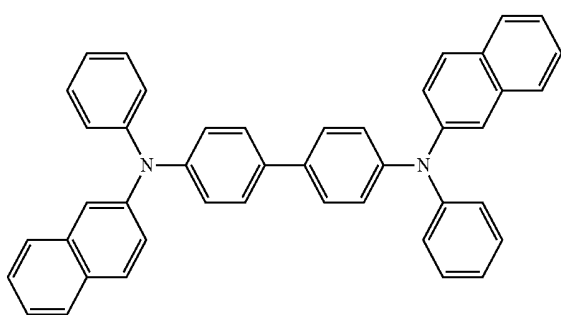
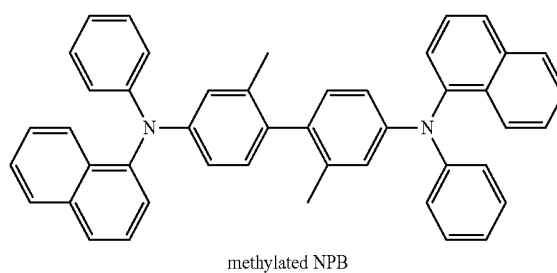
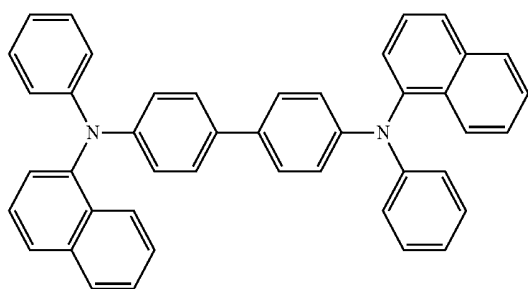
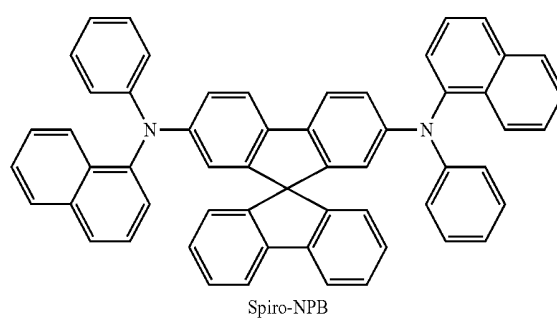
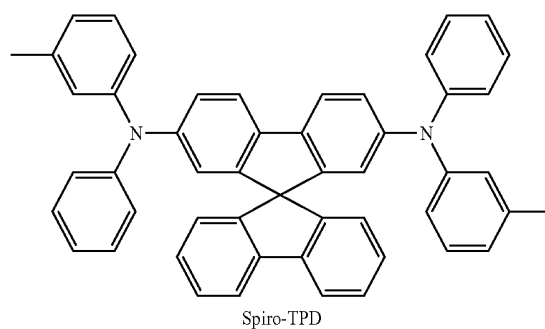
[0160] The hole transport region may include at least one selected from m-MTDATA, TDATA, 2-TNATA, NPB (NPD), 3-NPB, TPD, Spiro-TPD, Spiro-NPB, methylated-NPB, TAPC, HMTPD, 4,4',4''-tris(N-carbazolyl)triphenylamine (TCTA), polyaniline/dodecylbenzenesulfonic acid (PANI/DBSA), PEDOT/PSS (poly(3,4-ethylenedioxythiophene)/poly(4-styrenesulfonate)), polyaniline/camphor sulfonic acid (PANI/CSA), polyaniline/poly(4-styrenesulfonate) (PANI/PSS), a compound represented by Formula 201, and a compound represented by Formula 202:



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group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, a pyridinyl group, $-\text{Si}(\text{Q}_{31})(\text{Q}_{32})(\text{Q}_{33})$, and $-\text{N}(\text{Q}_{31})(\text{Q}_{32})$; and

[0171] Q_{31} to Q_{33} may each independently be selected from a C_1 - C_{10} alkyl group, a C_1 - C_{10} alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group.

[0172] In one or more embodiments, xa1 to xa4 may each independently be 0, 1, or 2.

[0173] In one or more embodiments, xa5 may be 1, 2, 3, or 4.

[0174] In one or more embodiments, R_{201} to R_{204} and Q_{201} may each independently be selected from a phenyl group, a biphenyl group, a terphenyl 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-bifluorenyl 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 hexacacenyl group, a pentacacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, and a pyridinyl group; and

[0175] a phenyl group, a biphenyl group, a terphenyl 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-bifluorenyl 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 hexacacenyl group, a pentacacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, and a pyridinyl group, each substituted with at least one selected from deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C_1 - C_{20} alkyl group, a 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 biphenyl group, a terphenyl group, a phenyl group substituted with a C_1 - C_{10} alkyl group, a phenyl group substituted with $-\text{F}$, 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-bifluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phe-

naleny 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 hexacacenyl group, a pentacacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, a pyridinyl group, $-\text{Si}(\text{Q}_{31})(\text{Q}_{32})(\text{Q}_{33})$, and $-\text{N}(\text{Q}_{31})(\text{Q}_{32})$, and

[0176] Q_{31} to Q_{33} are the same as described above.

[0177] In one or more embodiments, at least one of R_{201} to R_{203} in Formula 201 may each independently be selected from:

[0178] a fluorenyl group, a spiro-bifluorenyl group, a carbazolyl group, a dibenzofuranyl group, and a dibenzothiophenyl group; and

[0179] a fluorenyl group, a spiro-bifluorenyl group, a carbazolyl group, a dibenzofuranyl group, and a dibenzothiophenyl group, each substituted with at least one selected from deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C_1 - C_{20} alkyl group, a 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 biphenyl group, a terphenyl group, a phenyl group substituted with a C_1 - C_{10} alkyl group, a phenyl group substituted with $-\text{F}$, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a carbazolyl group, a dibenzofuranyl group, and a dibenzothiophenyl group,

[0180] but embodiments of the present disclosure are not limited thereto.

[0181] In one or more embodiments, in Formula 202, i) R_{201} and R_{202} may be linked via a single bond, and/or ii) R_{23} and R_{24} may be linked via a single bond.

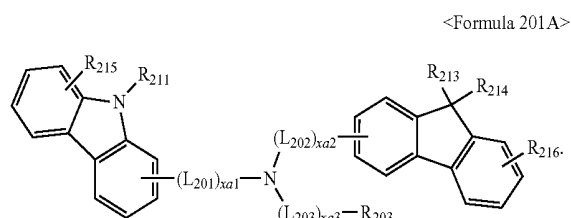
[0182] In one or more embodiments, at least one of R_{21} to R_{24} in Formula 202 may be selected from:

[0183] a carbazolyl group; and

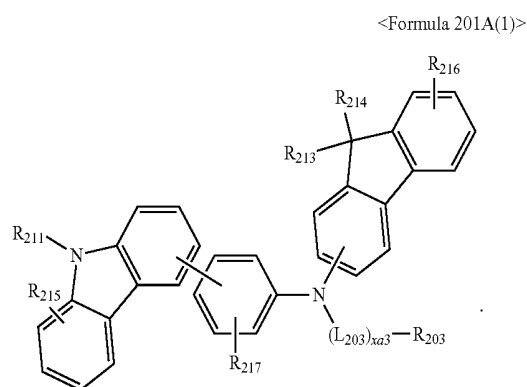
[0184] a carbazolyl group, substituted with at least one selected from deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl group, a biphenyl group, a terphenyl group, a phenyl group substituted with a C_1 - C_{10} alkyl group, a phenyl group substituted with $-\text{F}$, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a carbazolyl group, a dibenzofuranyl group, and a dibenzothiophenyl group,

[0185] but embodiments of the present disclosure are not limited thereto.

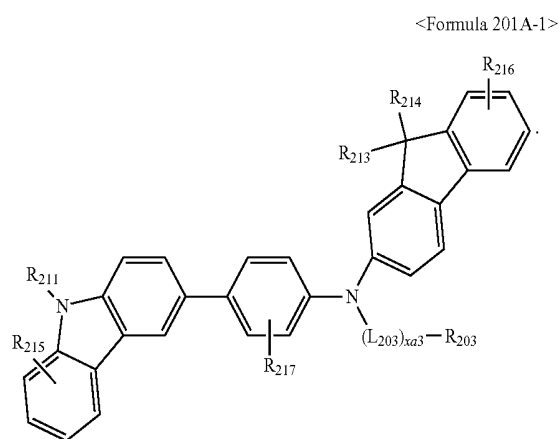
[0186] The compound represented by Formula 201 may be represented by Formula 201A:



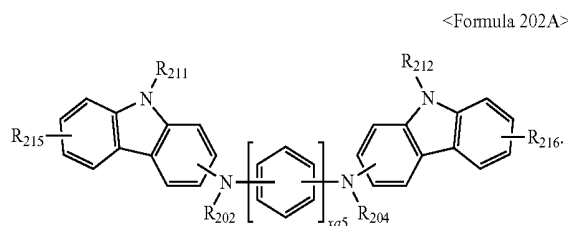
[0187] In one or more embodiments, the compound represented by Formula 201 may be represented by Formula 201A(1) below, but embodiments of the present disclosure are not limited thereto:



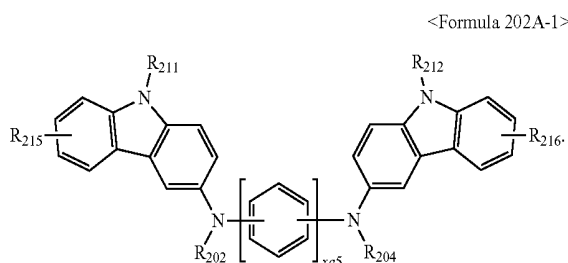
[0188] In one or more embodiments, the compound represented by Formula 201 may be represented by Formula 201A-1 below, but embodiments of the present disclosure are not limited thereto:



[0189] In one or more embodiments, the compound represented by Formula 202 may be represented by Formula 202A:



[0190] In one or more embodiments, the compound represented by Formula 202 may be represented by Formula 202A-1:



[0191] In Formulae 201A, 201A(1), 201A-1, 202A, and 202A-1,

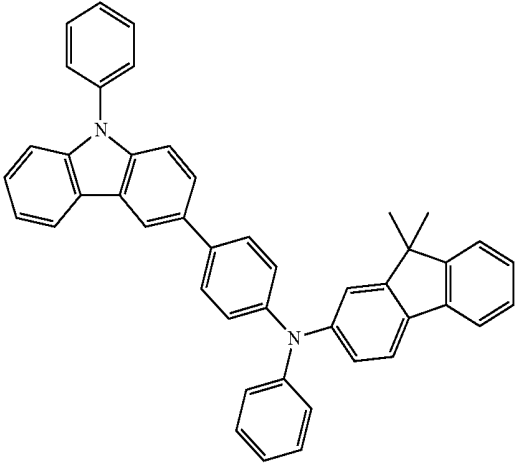
[0192] L_{201} to L_{203} , $xa1$ to $xa3$, $xa5$, and R_{202} to R_{204} are the same as described above, and

[0193] R_{211} and R_{212} may be understood by referring to the description provided herein in connection with R_{203} .

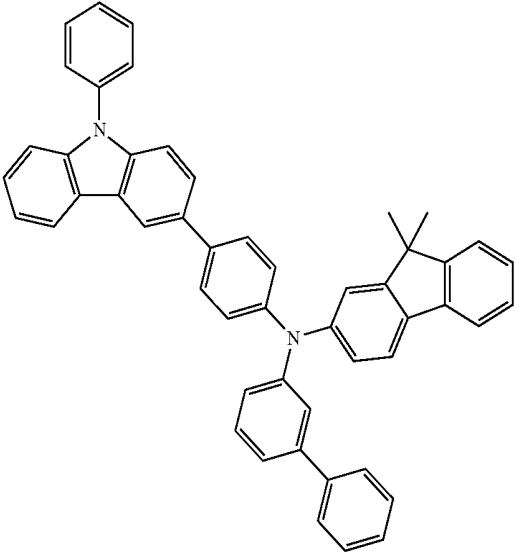
[0194] R_{213} to R_{217} may each independently be selected from hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl group, a biphenyl group, a terphenyl group, a phenyl group substituted with a C_1 - C_{10} alkyl group, a phenyl group substituted with —F, 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-bifluorenyl 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 thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, and a pyridinyl group.

[0195] The hole transport region may include at least one compound selected from Compounds HT1 to HT39, but embodiments of the present disclosure are not limited thereto:

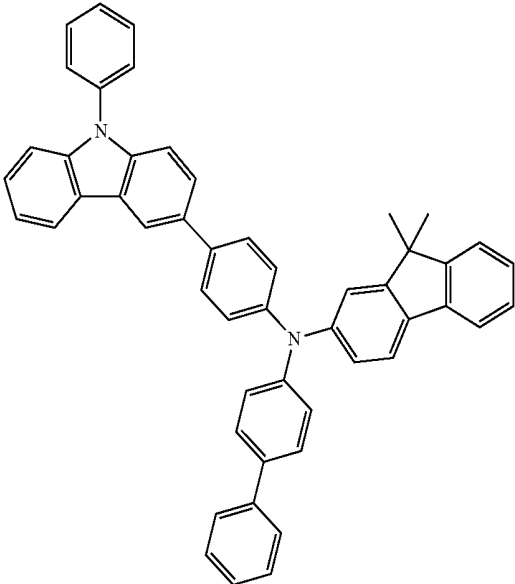
HT1



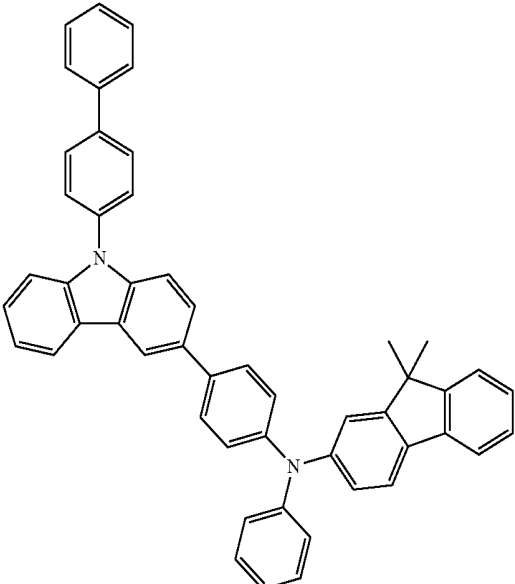
HT2



HT3

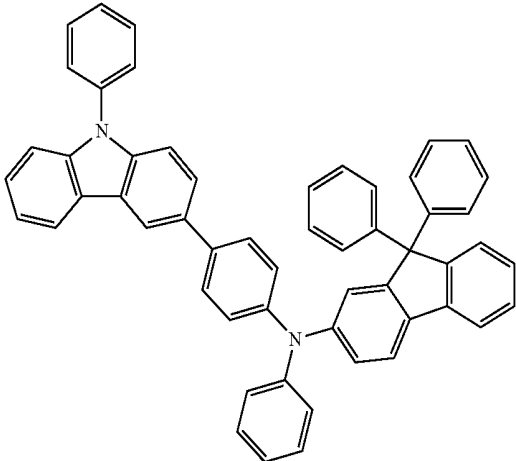


HT4

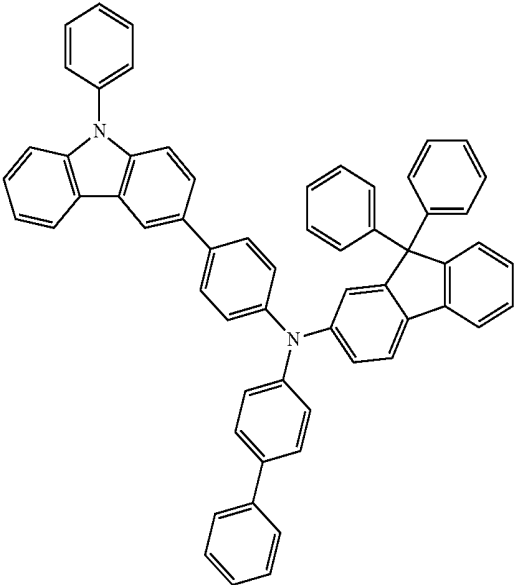


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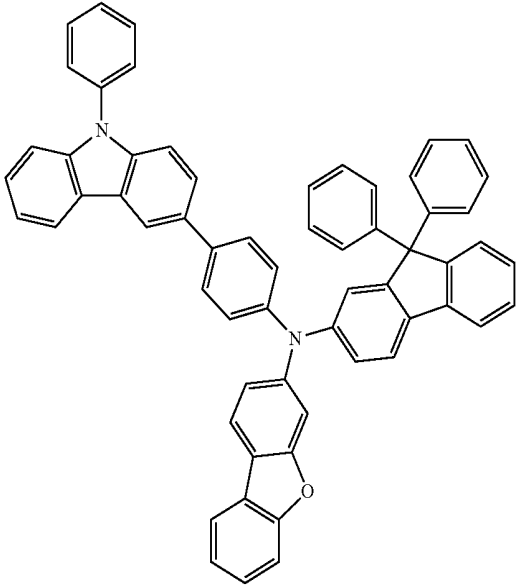
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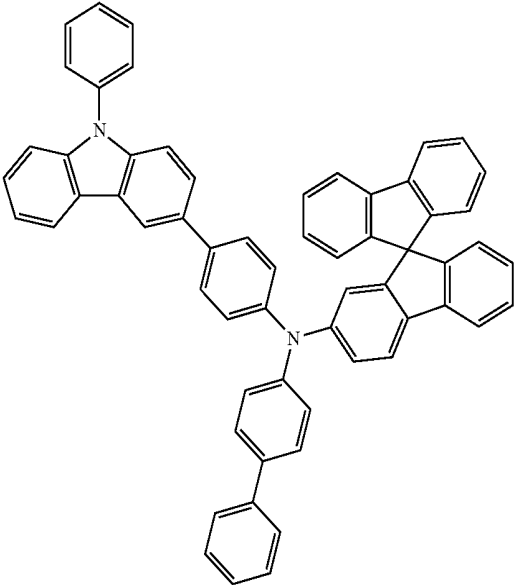
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HT7

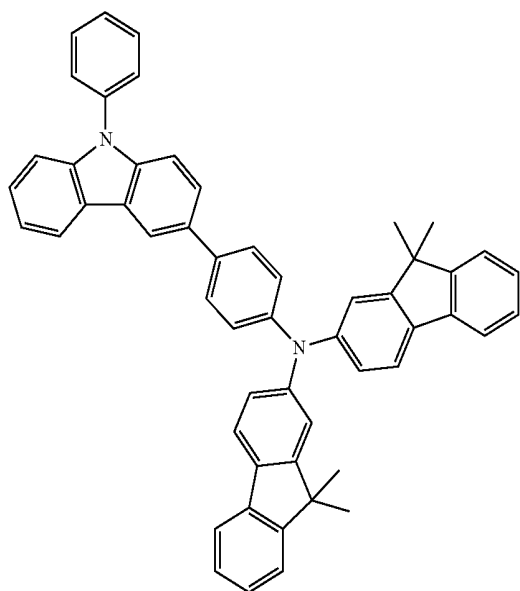


HT8

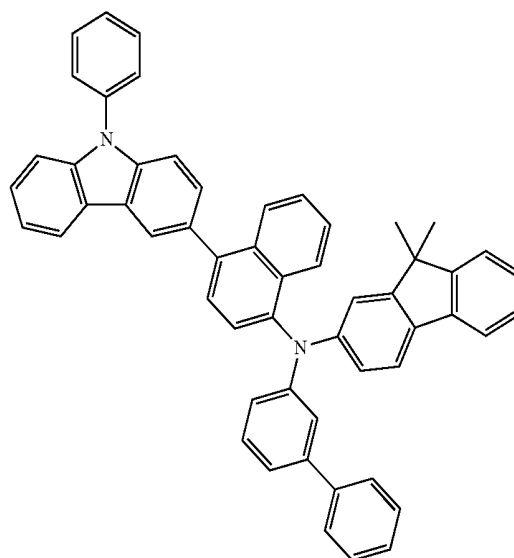


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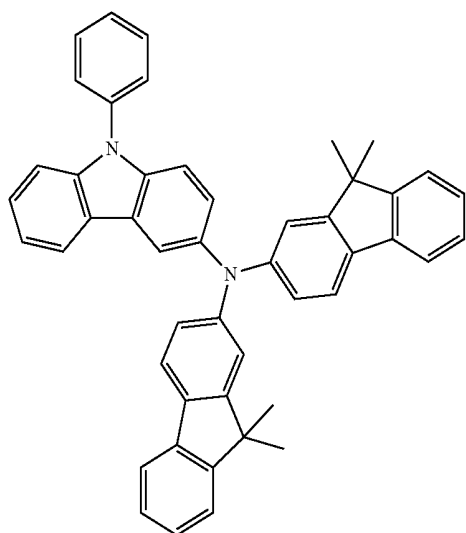
HT9



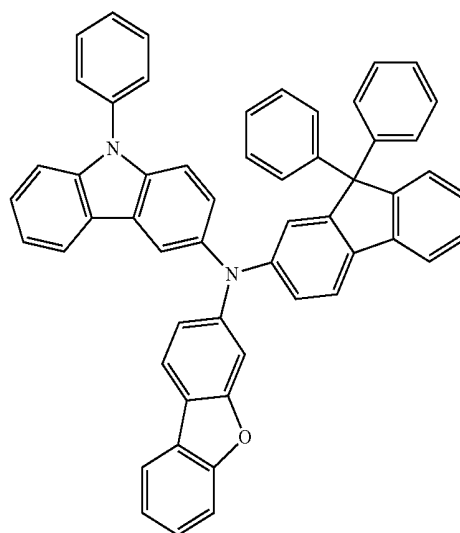
HT10



HT11

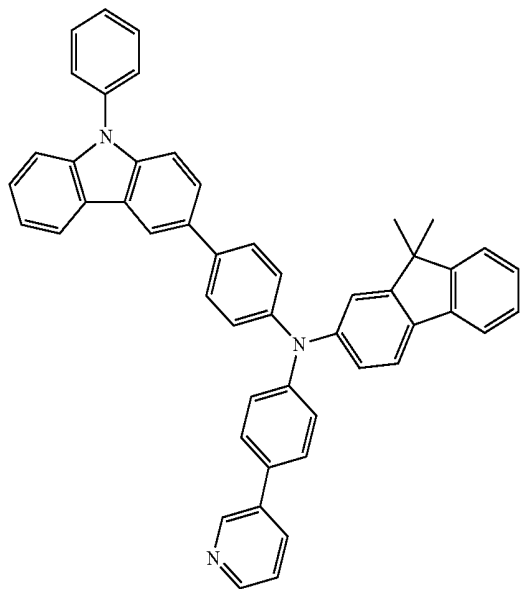


HT12

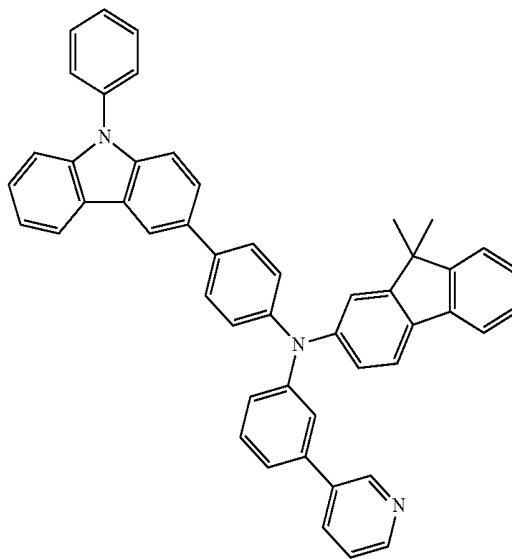


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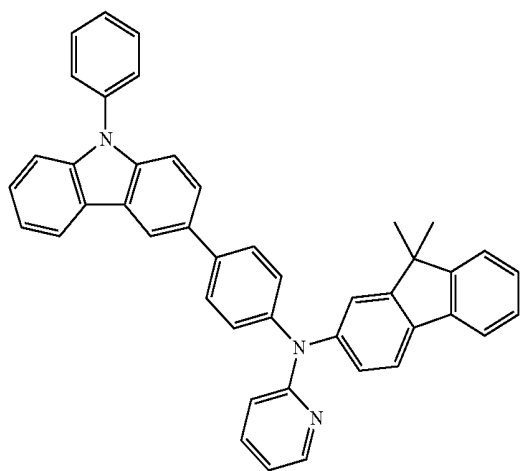
HT13



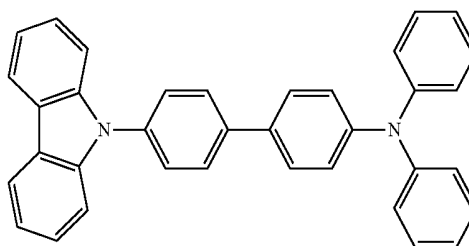
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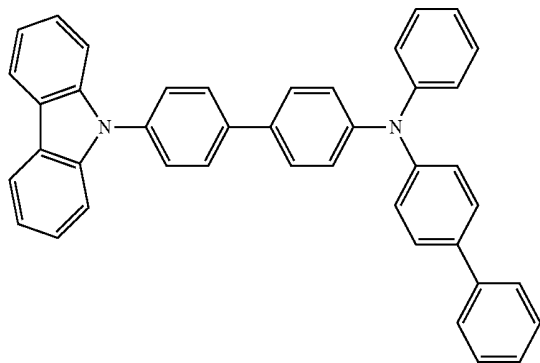
HT15



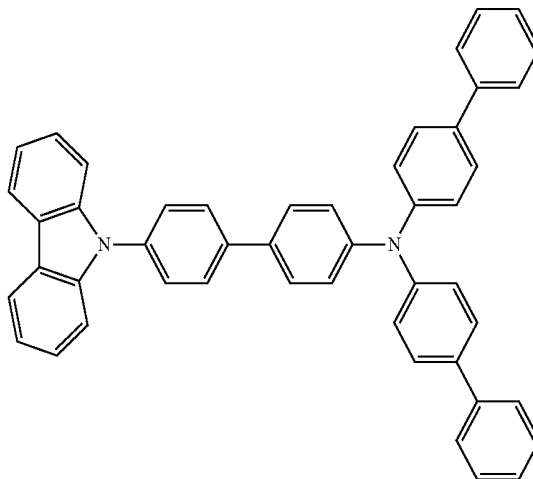
HT16



HT17



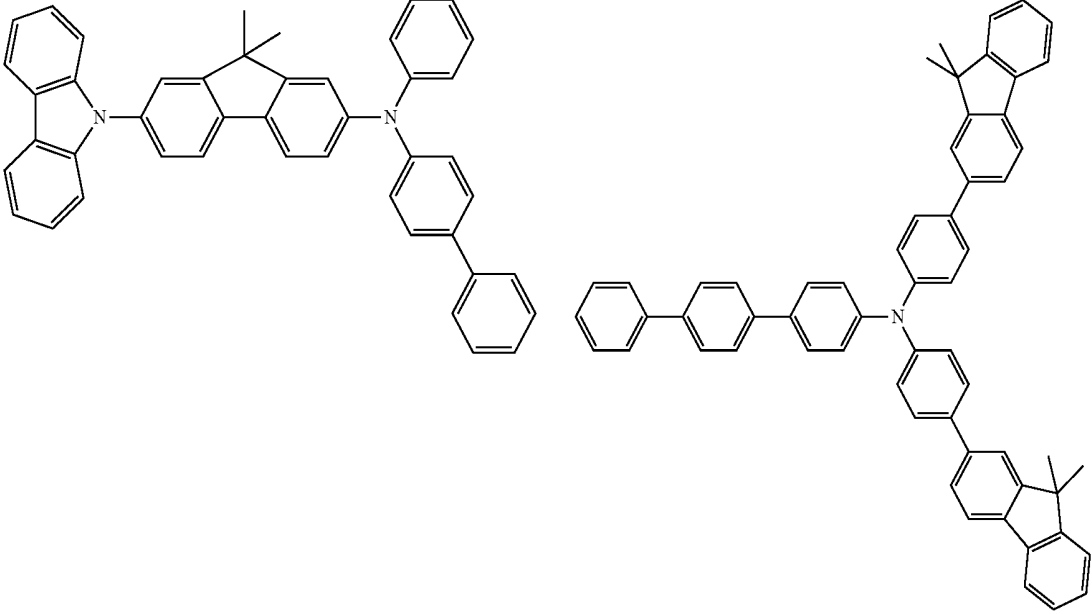
HT18



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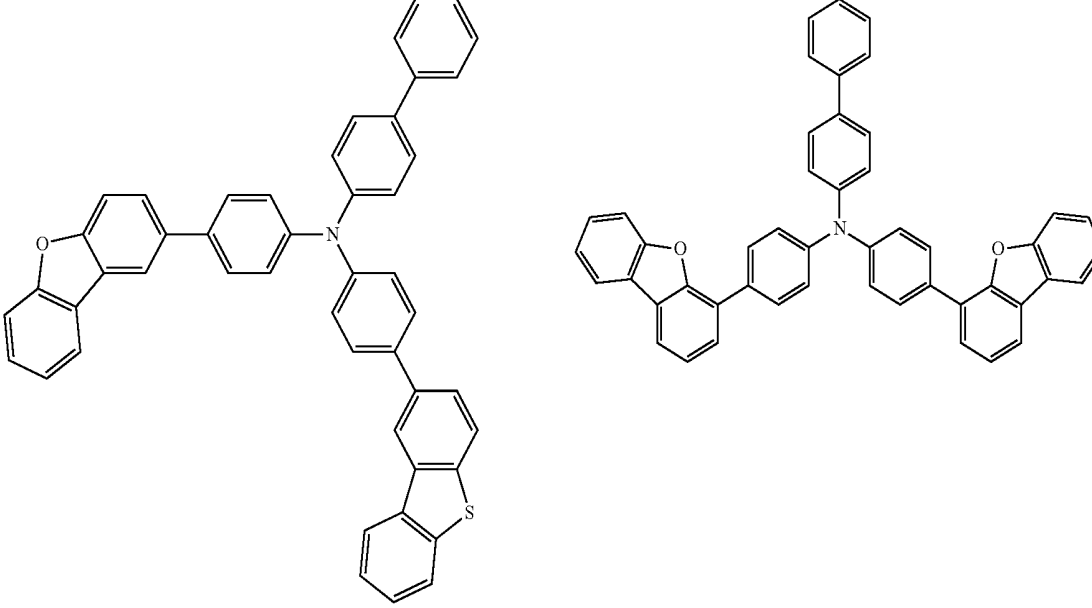
HT19

HT20



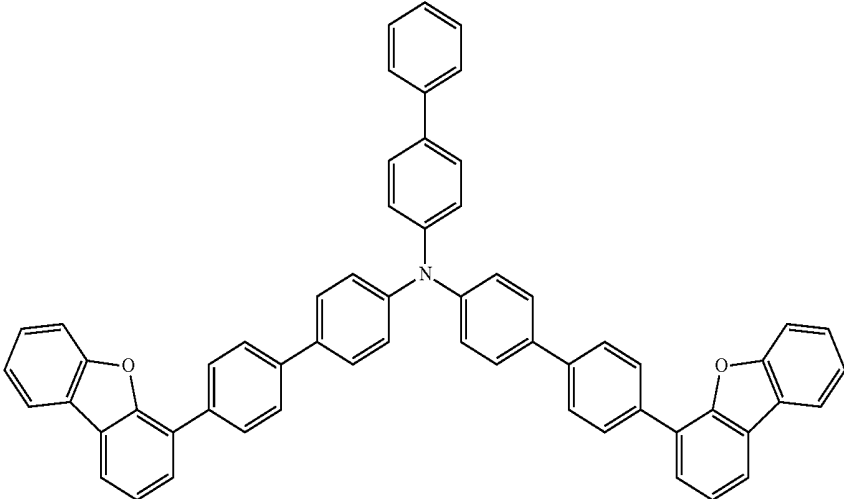
HT21

HT22

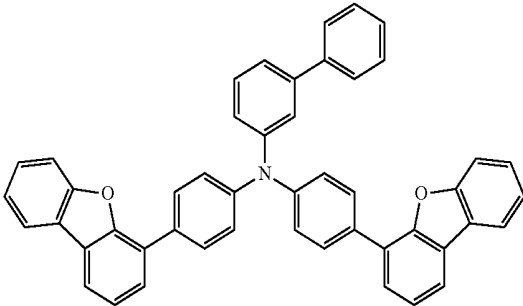


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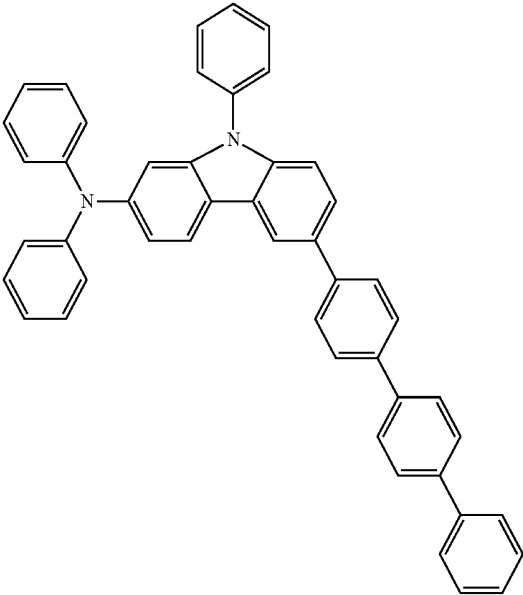
HT23



HT24

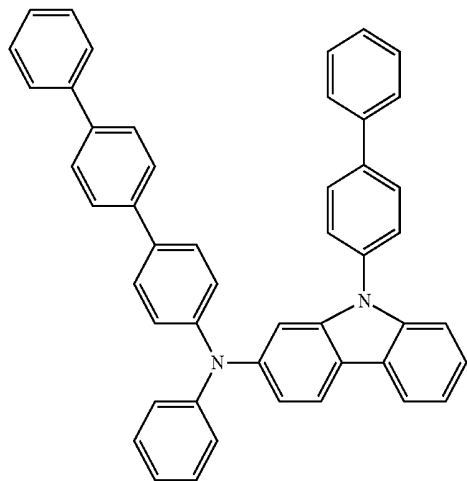


HT25

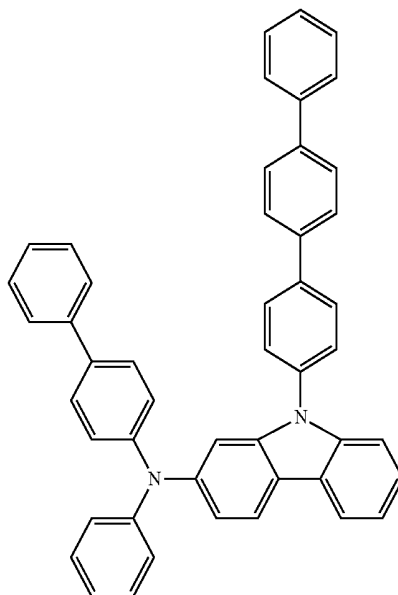


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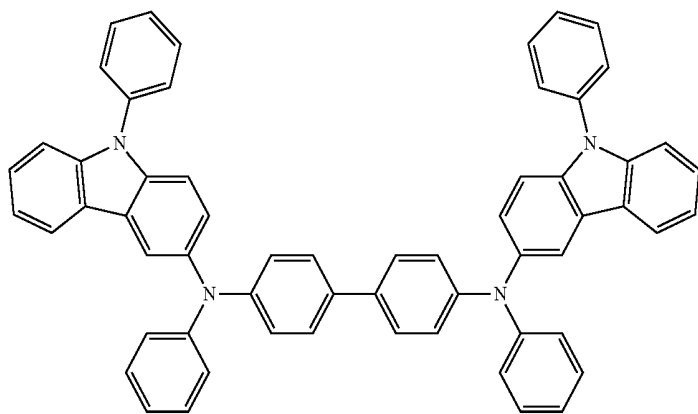
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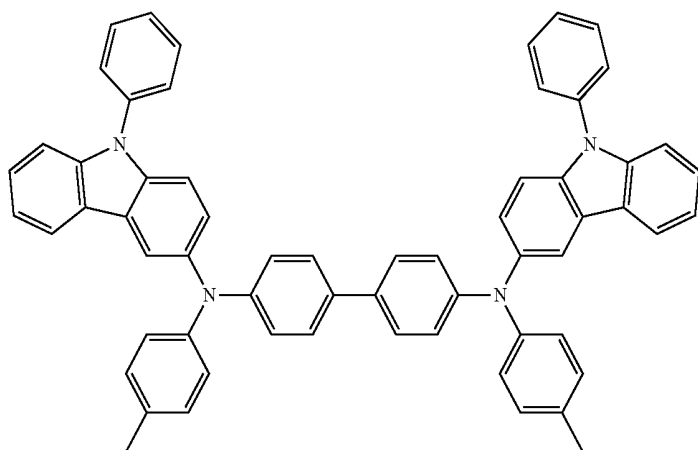
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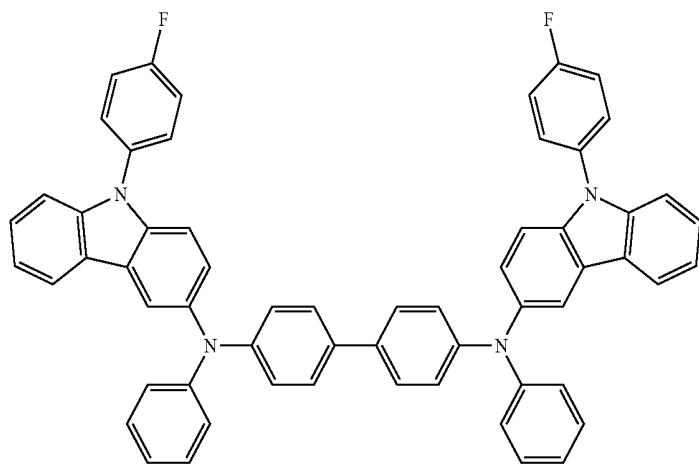
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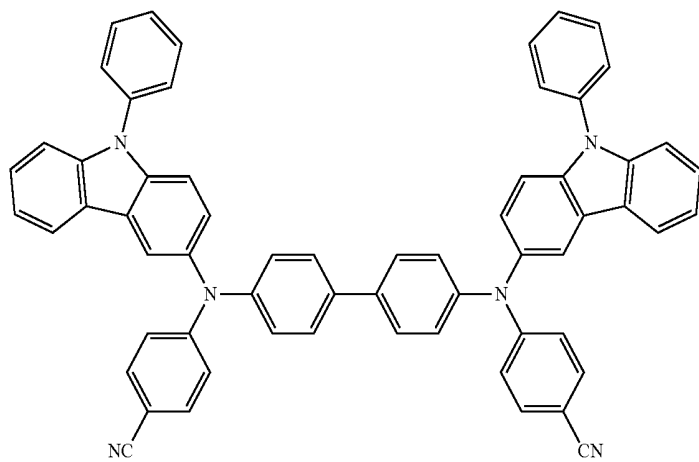
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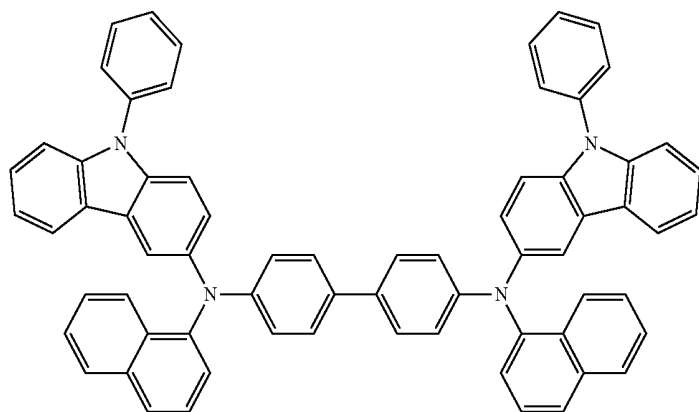
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HT30



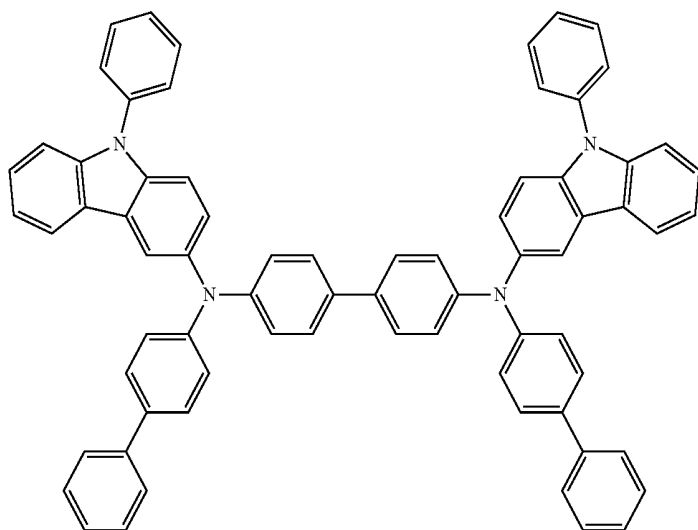
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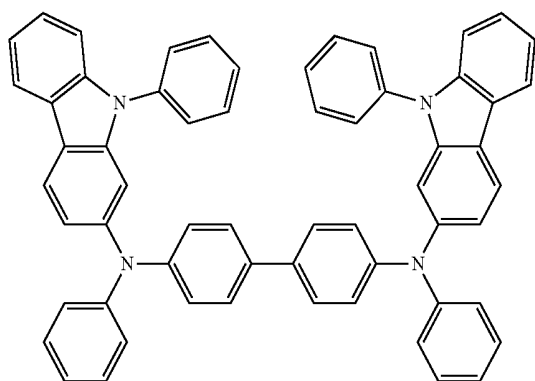
HT32

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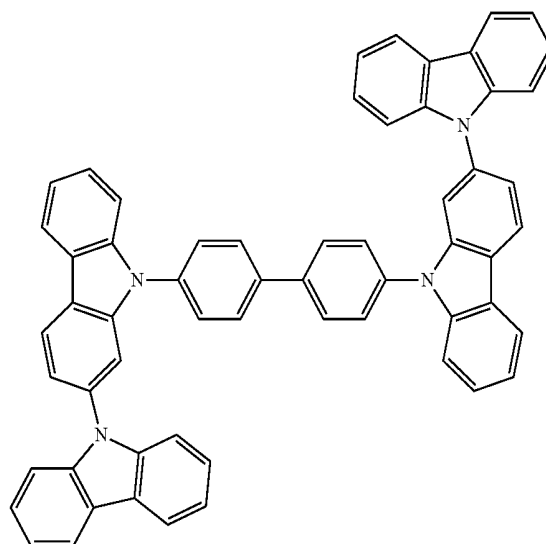
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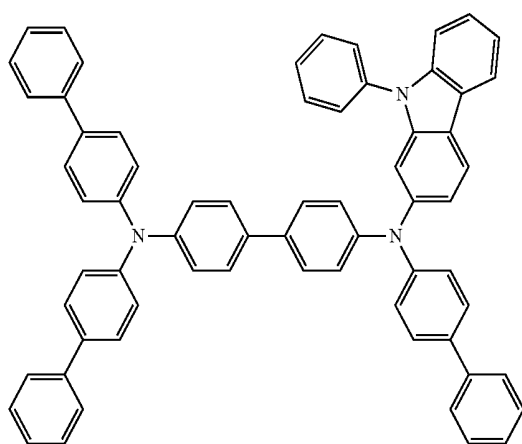
HT34



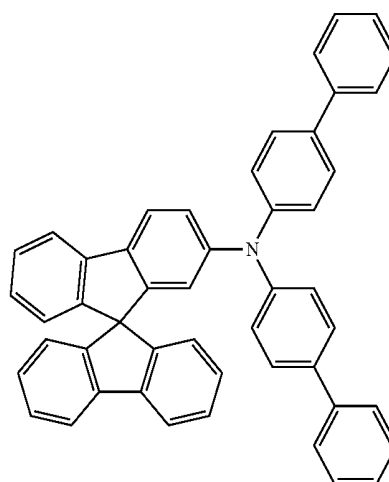
HT35



HT36

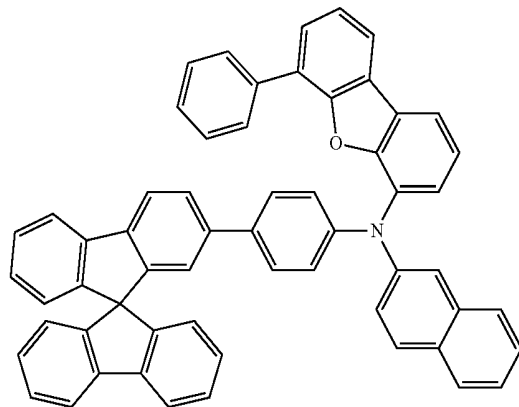
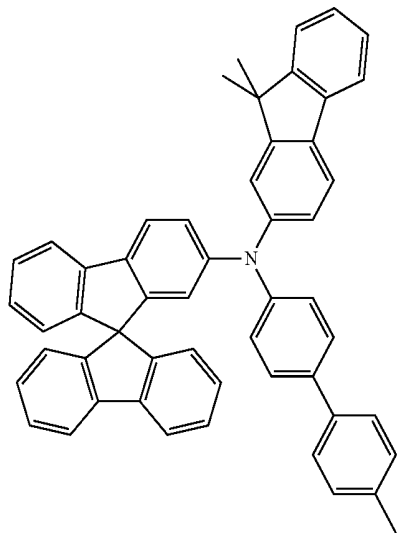


HT37



-continued
HT38

HT39



[0196] A thickness of the hole transport region may be in a range of about 100 Å to about 10,000 Å, for example, about 100 Å to about 1,000 Å. When the hole transport region includes at least one of a hole injection layer and a hole transport layer, the thickness of the hole injection layer may be in a range of about 100 Å to about 9,000 Å, and for example, about 100 Å to about 1,000 Å, and the thickness of the hole transport layer may be in a range of about 50 Å to about 2,000 Å, and for example, about 100 Å to about 1,500 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.

[0197] The emission auxiliary layer may increase light-emission efficiency by compensating for an optical resonance distance according to the wavelength of light emitted by an emission layer, and the electron blocking layer may block the flow of electrons from an electron transport region. The emission auxiliary layer and the electron blocking layer may include the materials as described above.

[p-Dopant]

[0198] 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.

[0199] The charge-generation material may be, for example, a p-dopant.

[0200] In one or more embodiments, a lowest unoccupied molecular orbital (LUMO) of the p-dopant may be smaller than -3.5 eV.

[0201] The p-dopant may include at least one selected from a quinone derivative, a metal oxide, and a cyano group-containing compound, but embodiments of the present disclosure are not limited thereto.

[0202] For example, the p-dopant may include at least one selected from:

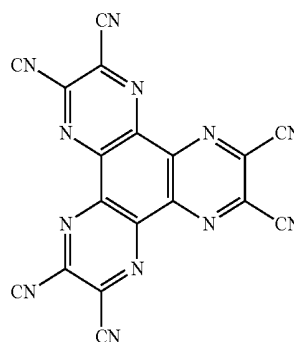
[0203] a quinone derivative, such as tetracyanoquinodimethane (TCNQ) and 2,3,5,6-tetrafluoro-7,7,8,8-tetracyanoquinodimethane (F4-TCNQ);

[0204] a metal oxide, such as tungsten oxide or molybdenum oxide;

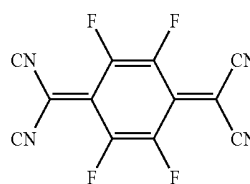
[0205] 1,4,5,8,9,11-hexaazatriphenylene-hexacarbonitrile (HAT-CN); and

[0206] a compound represented by Formula 221,

[0207] but embodiments of the present disclosure are not limited thereto:

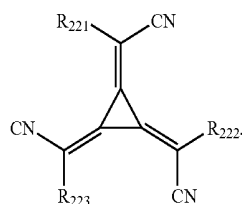


<HAT-CN>



<F4-TCNQ>

-continued



<Formula 221>

[0208] In Formula 221,

[0209] R_{221} to R_{223} may each independently 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_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, provided that at least one selected from R_{221} to R_{223} has at least one substituent selected from a cyano group, $-F$, $-Cl$, $-Br$, $-I$, a C_1 - C_{20} alkyl group substituted with $-F$, a C_1 - C_{20} alkyl group substituted with $-Cl$, a C_1 - C_{20} alkyl group substituted with $-Br$, and a C_1 - C_{20} alkyl group substituted with $-I$.

[Emission Layer in Organic Layer 150]

[0210] 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, according to a sub-pixel. In one or more embodiments, the emission layer may have a stacked structure of two or more layers selected from a red emission layer, a green emission layer, and a blue emission layer, in which the two or more layers contact each other or are separated from each other. In one or more embodiments, the emission layer may include two or more materials selected from a red light-emitting material, a green light-emitting material, and a blue light-emitting material, in which the two or more materials are mixed with each other in a single layer to emit white light.

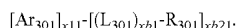
[0211] The emission layer may include a host and a dopant. The dopant may include at least one selected from a phosphorescent dopant and a fluorescent dopant.

[0212] An amount of the dopant in the emission layer may be, in general, in a range of about 0.01 parts by weight to about 15 parts by weight based on 100 parts by weight of the host, but embodiments of the present disclosure are not limited thereto.

[0213] A thickness of the emission layer may be in a range of about 100 Å to about 1,000 Å, for example, 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.

[Host in Emission Layer]

[0214] In one or more embodiments, the host may include a compound represented by Formula 301 below.



<Formula 301>

[0215] In Formula 301,

[0216] Ar_{301} may be a substituted or unsubstituted C_5 - C_{60} carbocyclic group or a substituted or unsubstituted C_1 - C_{60} heterocyclic group,

[0217] x_{b11} may be 1, 2, or 3,

[0218] L_{301} may be 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;

[0219] x_{b1} may be an integer selected from 0 to 5,

[0220] R_{301} may be selected from deuterium, $-F$, $-Cl$, $-Br$, $-I$, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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, $-Si(Q_{301})(Q_{302})(Q_{303})$, $-N(Q_{301})(Q_{302})$, $-B(Q_{301})(Q_{302})$, $-C(=O)(Q_{301})$, $-S(=O)_2(Q_{301})$, and $-P(=O)(Q_{301})(Q_{302})$.

[0221] x_{b21} may be an integer selected from 1 to 5, and

[0222] Q_{301} to Q_{303} may each independently be selected from a C_1 - C_{10} alkyl group, a C_1 - C_{10} alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group, but embodiments of the present disclosure are not limited thereto.

[0223] In one or more embodiments, Ar_{301} in Formula 301 may be selected from:

[0224] a naphthalene group, a fluorene group, a spiro-bifluorene 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, an indenoanthracene group, a dibenzofuran group, and a dibenzothiophene group; and

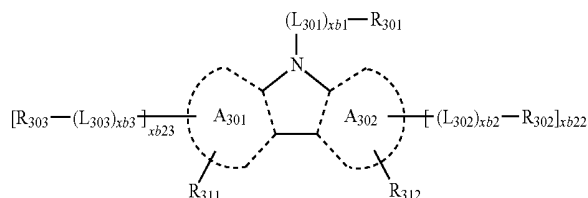
[0225] a naphthalene group, a fluorene group, a spiro-bifluorene group, a benzofluorene group, a dibenzofluorene group, a phenalene group, a phenanthrene group, an anthracene group, a fluoranthene group, a triphenylene group, a pyrene group, a chrysene group, a naphthacene group, a picene group, a perylene group, a pentaphene group, an indenoanthracene group, a dibenzofuran group, and a dibenzothiophene group,

each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, —Si(Q₃₁)(Q₃₂)(Q₃₃), —N(Q₃₁)(Q₃₂), —B(Q₃₁)(Q₃₂), —C(=O)(Q₃₁), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂); and

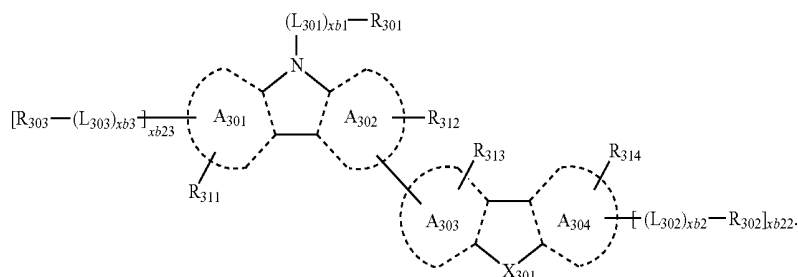
[0226] Q₃₁ to Q₃₃ may each independently be selected from a C₁-C₁₀ alkyl group, a C₁-C₁₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group. However, embodiments of the present disclosure are not limited thereto.

[0227] When xb11 in Formula 301 is two or more, two or more Ar₃₀₁(s) may be linked via a single bond.

[0228] In one or more embodiments, the compound represented by Formula 301 may be represented by Formula 301-1 or 301-2:



<Formula 301-1>



<Formula 301-2>

[0229] In Formulae 301-1 and 301-2,

[0230] A₃₀₁ to A₃₀₄ may each independently be selected from a benzene group, a naphthalene group, a phenanthrene group, a fluoranthene group, a triphenylene group, a pyrene group, a chrysene group, a pyridine group, a pyrimidine group, an indene group, a fluorene group, a spiro-bifluorene group, a benzofluorene group, a dibenzofluorene group, an indole group, a carbazole group, a benzocarbazole group, a dibenzocarbazole group, a furan group, a benzofuran group, a dibenzofuran group, a naphthofuran group, a benzonaphthofuran group, a dinaphthofuran group, a thiophene group, a benzothiophene group, a dibenzothiophene group, a naphthothiophene group, a benzonaphthothiophene group, and a dinaphthothiophene group,

[0231] X₃₀₁ may be O, S, or N-[(L₃₀₄)_{xb4}-R₃₀₄],

[0232] R₃₁₁ to R₃₁₄ may each independently be selected from hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, a naphthyl

group —Si(Q₃₁)(Q₃₂)(Q₃₃), —N(Q₃₁)(Q₃₂), —B(Q₃₁)(Q₃₂), —C(=O)(Q₃₁), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂),

[0233] xb22 and xb23 may each independently be 0, 1, or 2,

[0234] L₃₀₁, xb1, R₃₀₁, and Q₃₁ to Q₃₃ are the same as described above,

[0235] L₃₀₂ to L₃₀₄ may each independently be the same as described in connection with L₃₀₁,

[0236] xb2 to xb4 may each independently be the same as described in connection with xb1, and

[0237] R₃₀₂ to R₃₀₄ may each independently be the same as described in connection with R₃₀₁.

[0238] For example, L₃₀₁ to L₃₀₄ in Formulae 301, 301-1, and 301-2 may each independently be selected from:

[0239] a phenylene group, a naphthylene group, a fluorenylene group, a spiro-bifluorenylene group, a benzofluorenylene group, a dibenzofluorenylene group, a

phenanthrenylene group, an anthracenylene group, a fluoranthenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a perylenylene group, a pentaphenylenylene group, a hexacenylenylene group, a pentacenylenylene group, a thiophenylenylene group, a furanylenylene group, a carbazolylenylene group, an indolylenylene group, an isoindolylenylene group, a benzofuranylenylene group, a benzothiophenylenylene group, a dibenzofuranylenylene group, a dibenzothiophenylenylene group, a benzocarbazolylenylene group, a dibenzocarbazolylenylene group, a dibenzosilolylenylene group, a pyridinylenylene group, an imidazolylenylene group, a pyrazolylenylene group, a thiazolylenylene group, an isothiazolylenylene group, an oxazolylenylene group, an isoxazolylenylene group, a thiadiazolylenylene group, an oxadiazolylenylene group, a pyrazinylenylene group, a pyrimidinylenylene group, a pyridazinylenylene group, a triazinylenylene group, a quinolininylenylene group, an isoquinolininylenylene group, a benzoquinolininylenylene group, a phthalazinylenylene group, a naphthyridinylenylene group, a quinoxalinylenylene group, a quinazolinylenylene group, a cinolininylenylene group, a phenanthridinylenylene group, an acridinylenylene group, a phenanthrolinylenylene group, a phenazinylenylene group, a benzimidazolylenylene group, an isobenzothiazolylenylene group, a benzoxazolylenylene group,

an isobenzoxazolylene group, a triazolylene group, a tetrazolylene group, an imidazopyridinylene group, an imidazopyrimidinylene group, and an azacarbazolylene group; and

[0240] a phenylene group, a naphthylene group, a fluorenylene group, a spiro-bifluorenylene group, a benzo-fluorenylene group, a dibenzofluorenylene group, a phenanthrenylene group, an anthracenylene group, a fluoranthenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a perylenylene group, a pentaphenylene group, a hexacenylenylene group, a pentacenylenylene group, a thiophenylene group, a furanylene group, a carbazolylene group, an indolylene group, an isoindolylene group, a benzofuranylene group, a benzothiophenylene group, a dibenzofuranylene group, a dibenzothiophenylene group, a benzocarbazolylene group, a dibenzocarbazolylene group, a dibenzosilolylene group, a pyridinylene group, an imidazolylene group, a pyrazolylene group, a thiazolylene group, an isothiazolylene group, an oxazolylene group, an isoxazolylene group, a thiadiazolylene group, an oxadiazolylene group, a pyrazinylene group, a pyrimidinylene group, a pyridazinylene group, a triazinylene group, a quinolinylene group, an isoquinolinylene group, a benzoquinolinylene group, a phthalazinylene group, a naphthyridinylene group, a quinoxalinylene group, a quinazolinylene group, a cinnolinylene group, a phenanthridinylene group, an acridinylene group, a phenanthrolinylene group, a phenazinylene group, a benzimidazolylene group, an isobenzothiazolylene group, a benzoxazolylene group, an isobenzoxazolylene group, a triazolylene group, a tetrazolylene group, an imidazopyridinylene group, an imidazopyrimidinylene group, and an azacarbazolylene group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, a pyridinyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a thiadiazolyl group, an oxadiazolyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a triazinyl 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 phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzimidazolyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, a tetrazolyl group, an imida-

zopyridinyl group, an imidazopyrimidinyl group, an azacarbazolyl group, —Si(Q₃₁)(Q₃₂)(Q₃₃), —N(Q₃₁)(Q₃₂), —B(Q₃₁)(Q₃₂), —C(=O)(Q₃₁), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂), and

[0241] Q₃₁ to Q₃₃ are the same as described above.

[0242] In one or more embodiments, R₃₀₁ to R₃₀₄ in Formulae 301, 301-1, and 301-2 may each independently be selected from:

[0243] a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, a pyridinyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a thiadiazolyl group, an oxadiazolyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a triazinyl 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 phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzimidazolyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, a tetrazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and an azacarbazolyl group; and

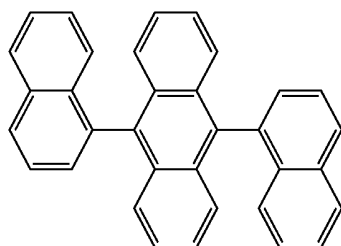
[0244] a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, a pyridinyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a thiadiazolyl group, an oxadiazolyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a triazinyl 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 phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzimidazolyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, a tetrazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group,

and an azacarbazolyl group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthrenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, a pyridinyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a thiadiazolyl group, an oxadiazolyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinoxalinyl group, a cinnolinyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzimidazolyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, a tetrazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, an azacarbazolyl group, $-\text{Si}(\text{Q}_{31})(\text{Q}_{32})(\text{Q}_{33})$, $-\text{N}(\text{Q}_{31})(\text{Q}_{32})$, $-\text{B}(\text{Q}_{31})(\text{Q}_{32})$, $-\text{C}(=\text{O})(\text{Q}_{31})$, $-\text{S}(=\text{O})_2(\text{Q}_{31})$, and $-\text{P}(=\text{O})(\text{Q}_{31})(\text{Q}_{32})$, and

[0245] Q_{31} and Q_{33} may be the same as described above.

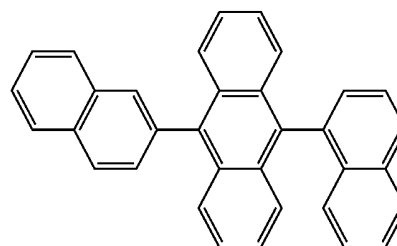
[0246] In one or more embodiments, the host may include an alkaline earth metal complex. For example, the host may be selected from a Be complex (for example, Compound H55), a Mg complex, and a Zn complex.

[0247] The host may include at least one selected from 9,10-di(2-naphthyl)anthracene (ADN), 2-methyl-9,10-bis(naphthalen-2-yl)anthracene (MADN), 9,10-di(2-naphthyl)-2-t-butyl-anthracene (TBADN), a 4,4'-bis(N-carbazolyl)-1,1'-biphenyl (CBP), 1,3-di-9-carbazolylbenzene (mCP), 1,3,5-tri(carbazol-9-yl)benzene (TCP), and Compounds H1 to H55, but embodiments of the present disclosure are not limited thereto:

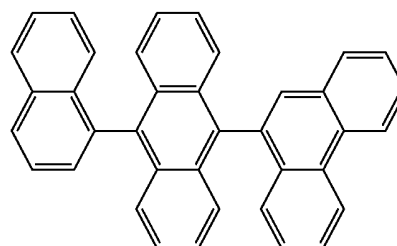


H1

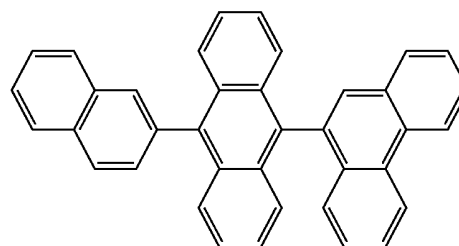
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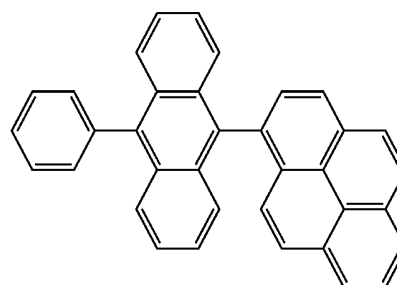
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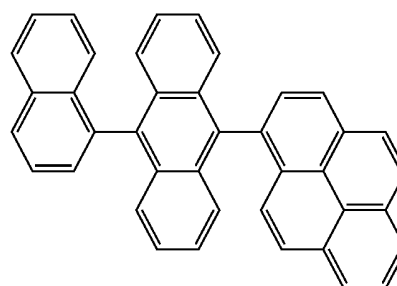
H3



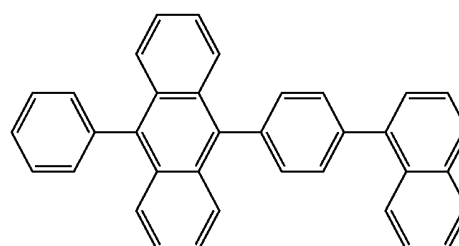
H4



H5



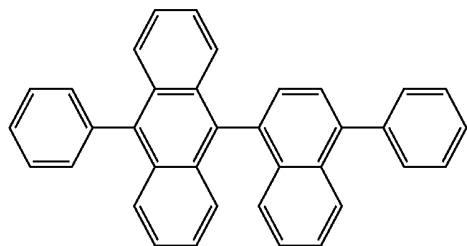
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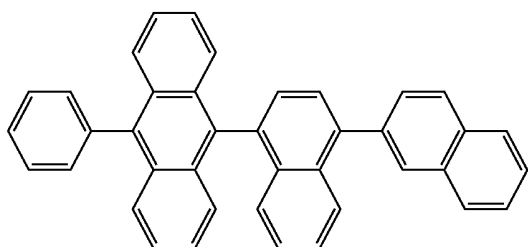
H7

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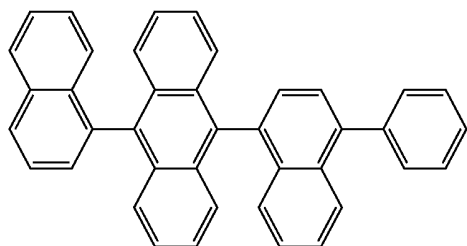
H8



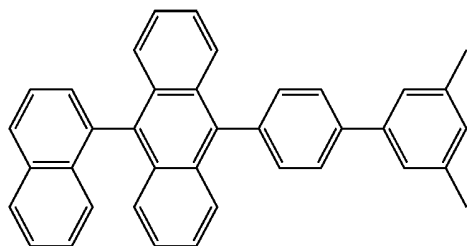
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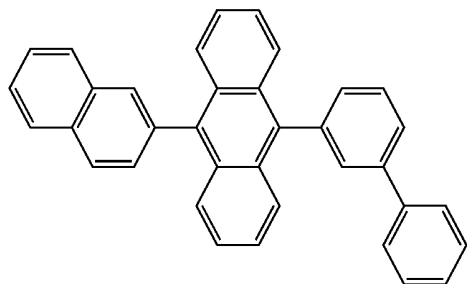
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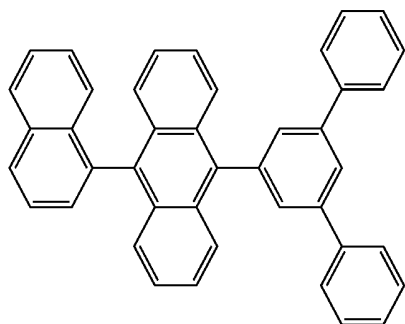
H11



H12

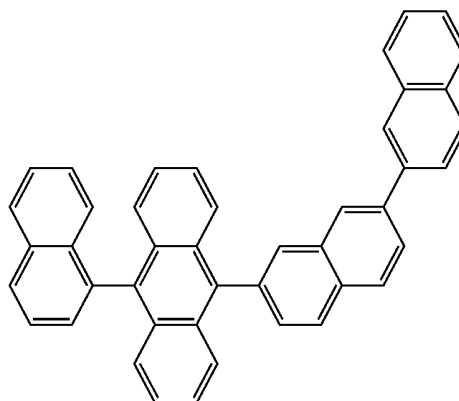


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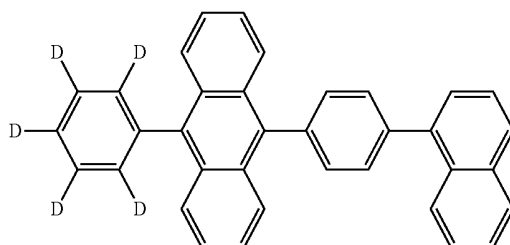


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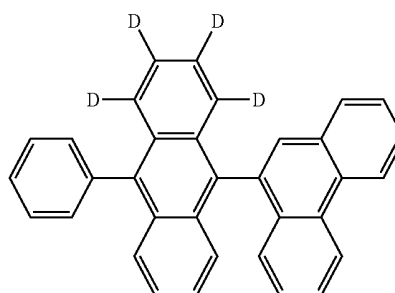
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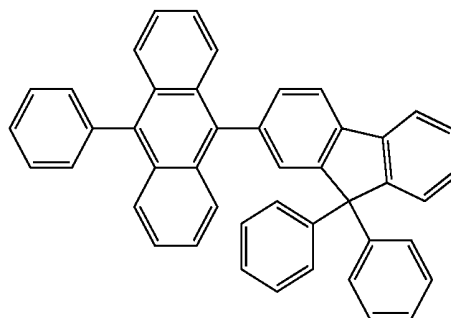
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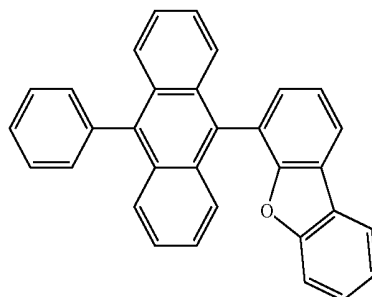
H16



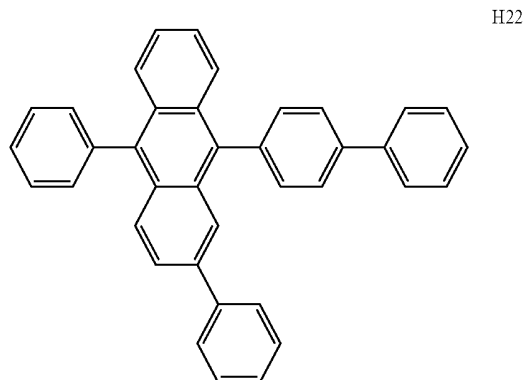
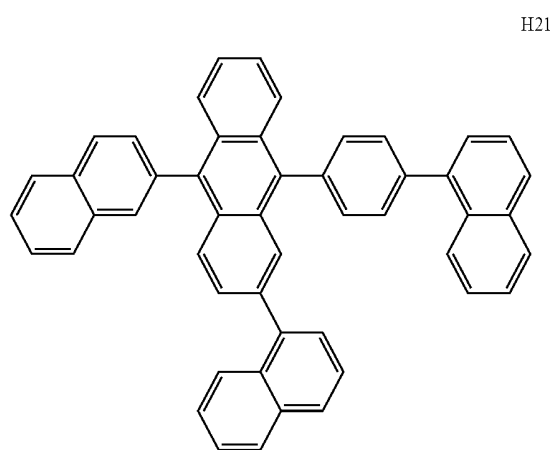
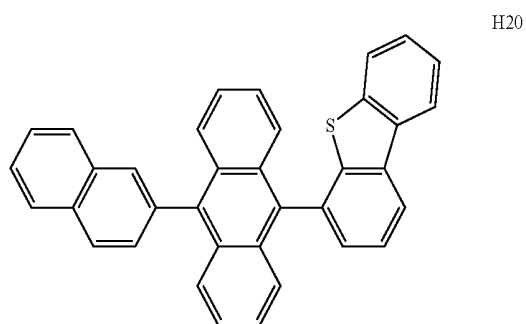
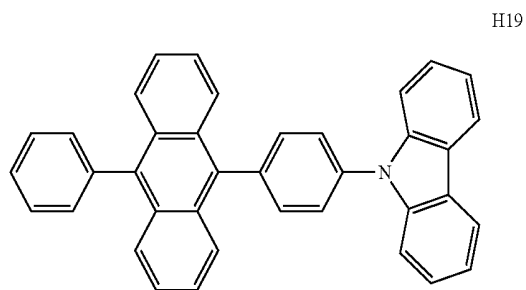
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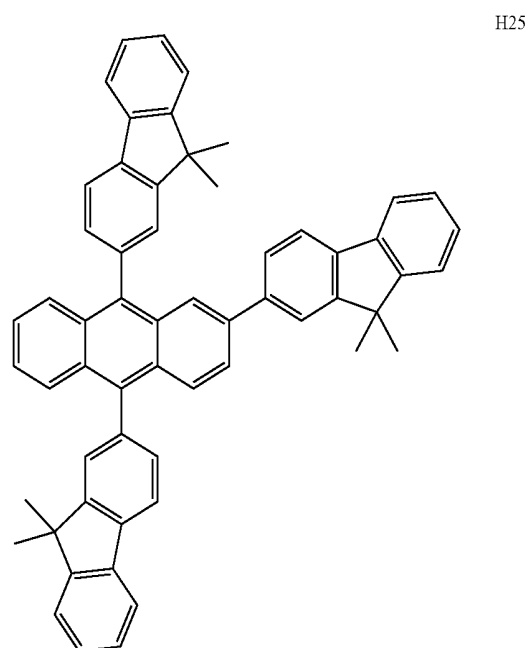
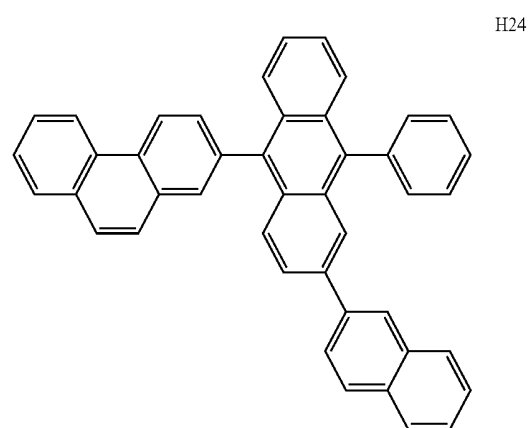
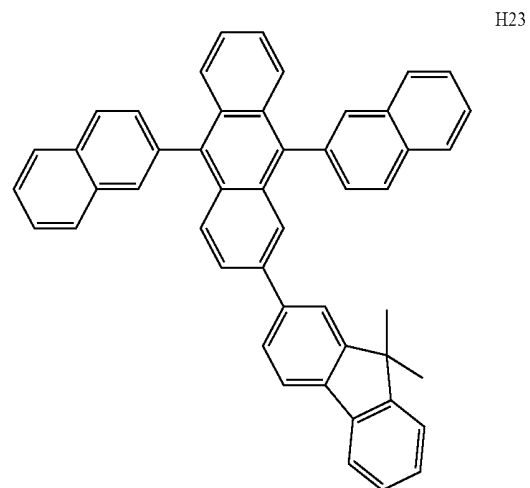
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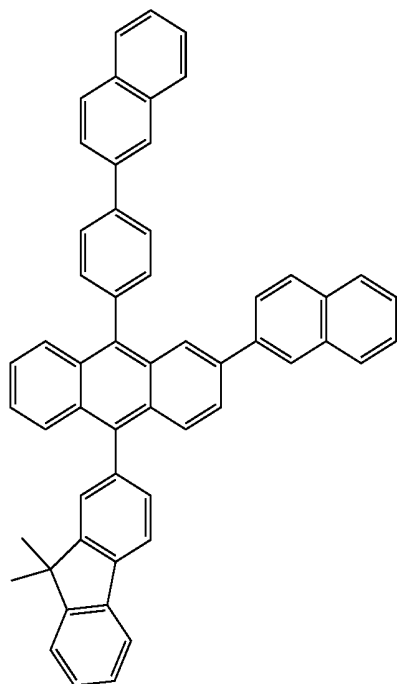
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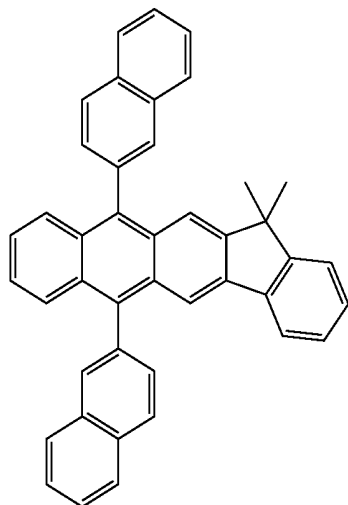
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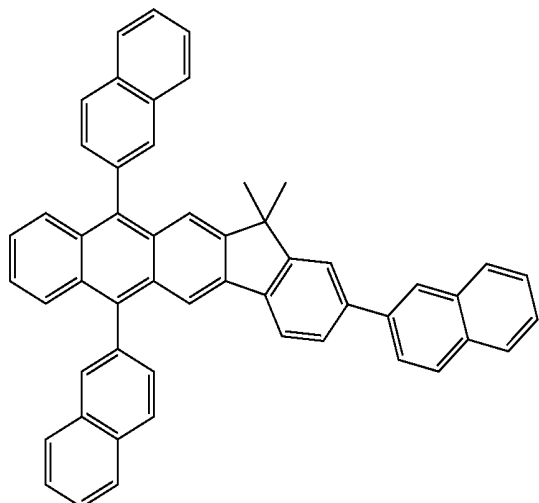
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H26

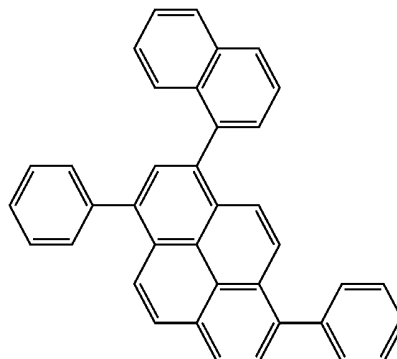


H27

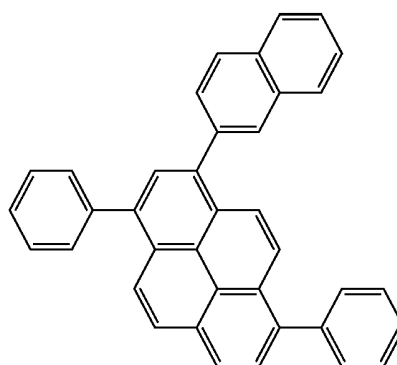


H28

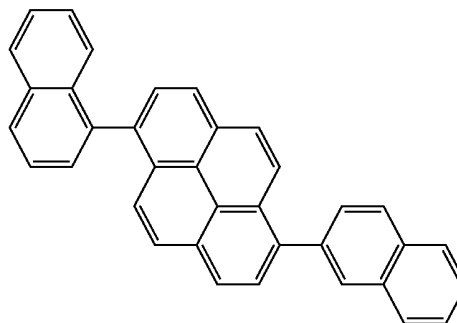
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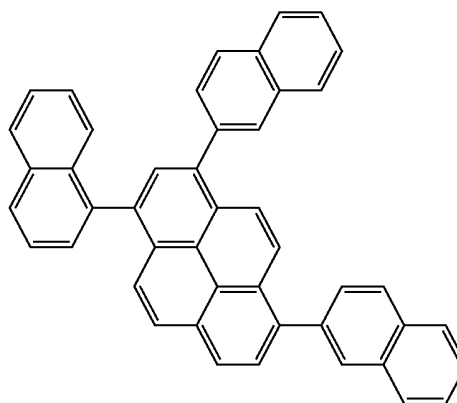
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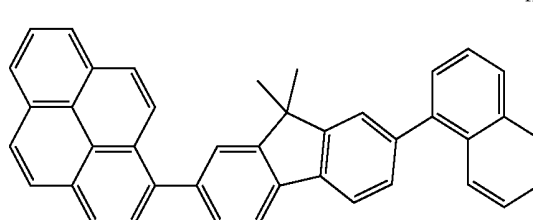
H30



H31



H32

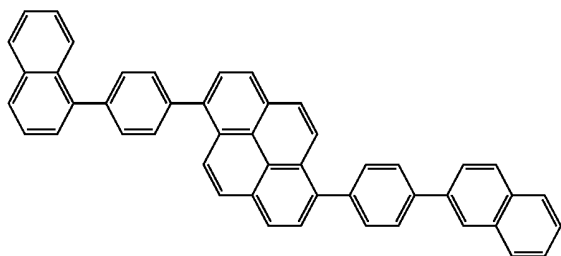


H33

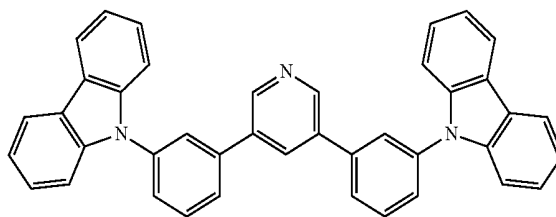
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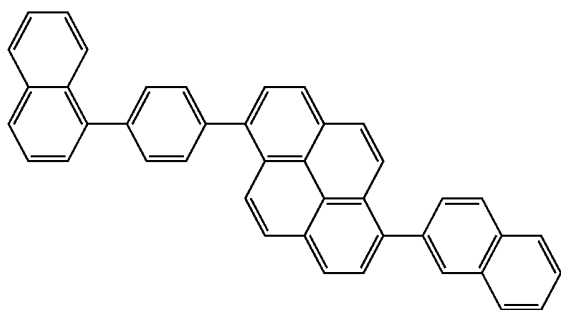
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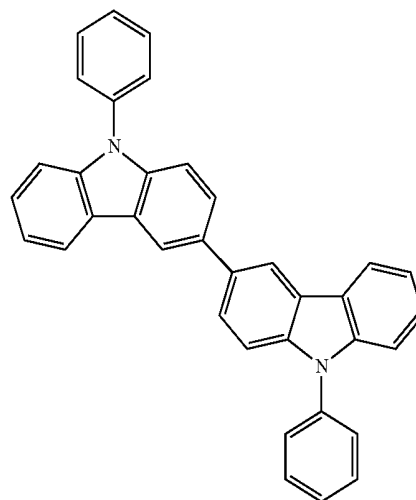
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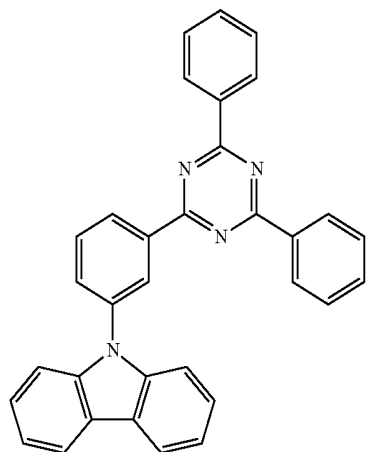
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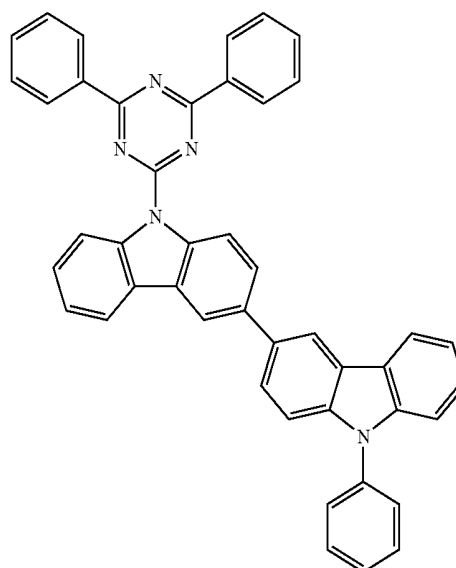
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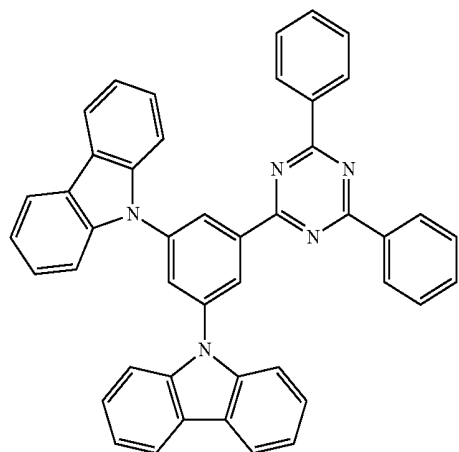
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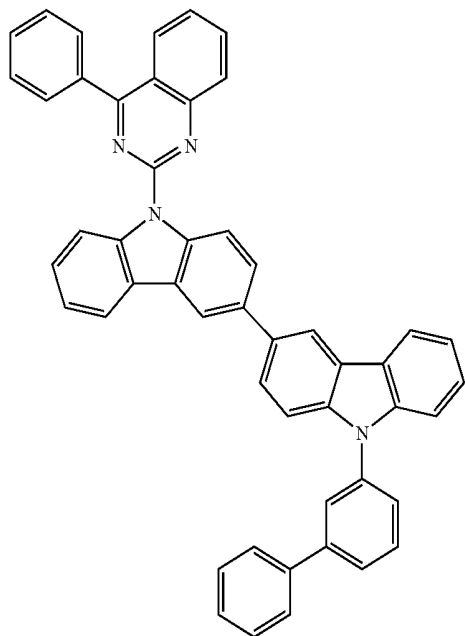
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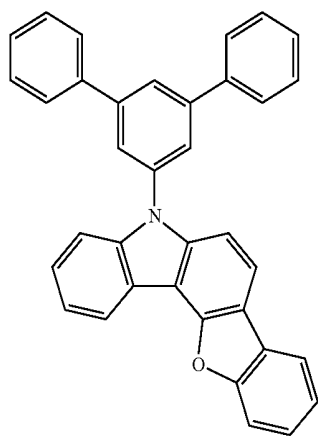
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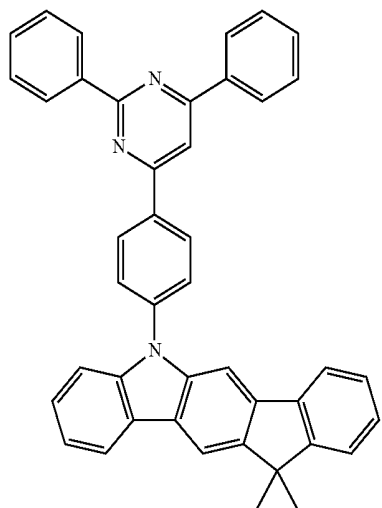
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H41

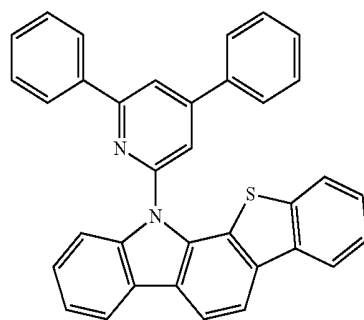


H42

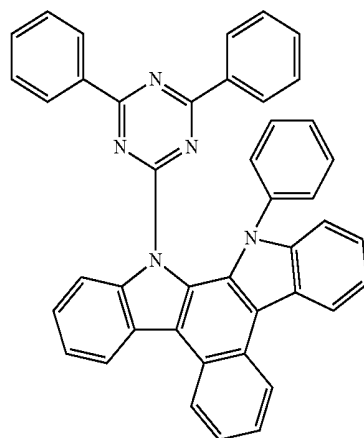


H43

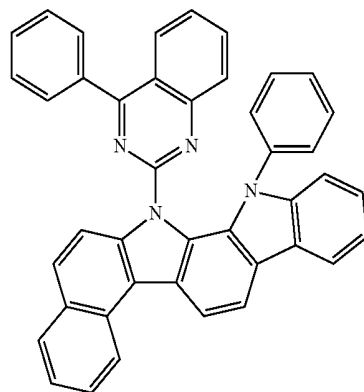
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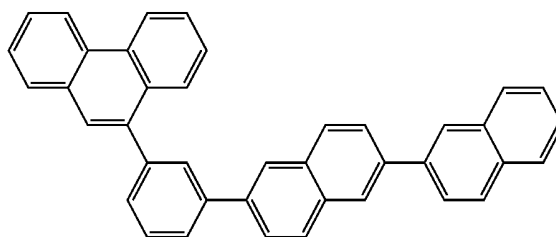
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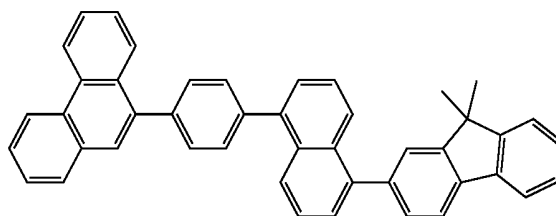
H45



H46

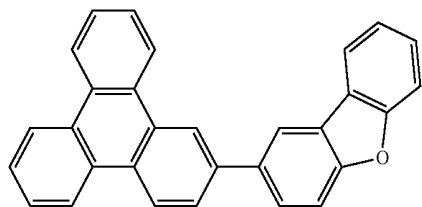


H47

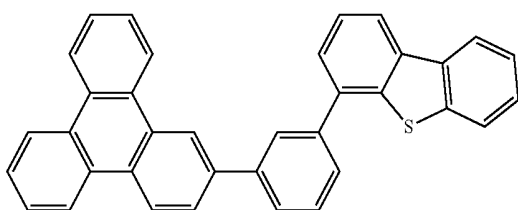


H48

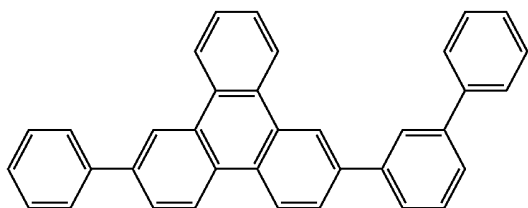
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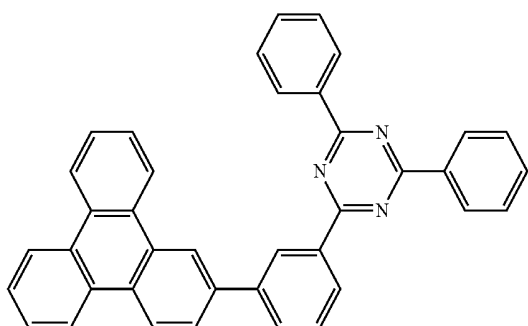
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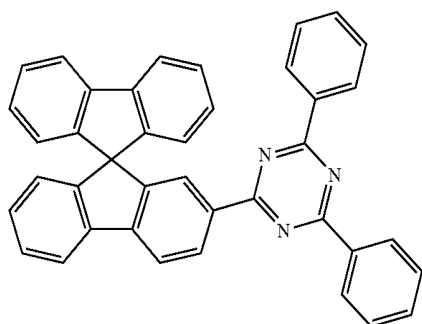
H50



H51

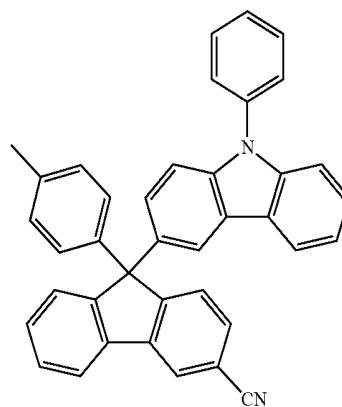


H52

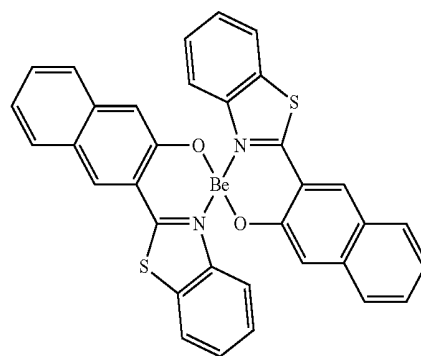


H53

-continued



H54



H55

[0248] In one or more embodiments, the host may include at least one selected from a silicon-containing compound (for example, BCPDS used in the following examples or the like) and a phosphine oxide-containing compound (for example, POPCPA used in the following examples or the like).

[0249] However, embodiments of the present disclosure are not limited thereto. In one or more embodiments, the host may include only one compound, or two or more different compounds (for example, a host used in the following examples includes BCPDS and POPCPA).

[Phosphorescent Dopant Included in Emission Layer in Organic Layer 150]

[0250] The phosphorescent dopant may include an organometallic compound represented by Formula 1 above.

[Electron Transport Region in Organic Layer 150]

[0251] The electron transport region may have i) a single-layered structure including a single layer including a single material, ii) a single-layered structure including a single layer including a plurality of different materials, or iii) a multi-layered structure having a plurality of layers including a plurality of different materials.

[0252] The electron transport region may include at least one selected from a buffer layer, a hole blocking layer, an electron control layer, an electron transport layer, and an electron injection layer, but embodiments of the present disclosure are not limited thereto.

[0253] For example, the electron transport region may have an electron transport layer/electron injection layer structure, a hole blocking layer/electron transport layer/

electron injection layer structure, an electron control layer/electron transport layer/electron injection layer structure, or a buffer layer/electron transport layer/electron injection layer structure, wherein for each structure, constituting layers are sequentially stacked from an emission layer. However, embodiments of the structure of the electron transport region are not limited thereto.

[0254] The electron transport region (for example, a buffer layer, a hole blocking layer, an electron control layer, or an electron transport layer in the electron transport region) may include a metal-free compound containing at least one π electron-depleted nitrogen-containing ring.

[0255] The “ π electron-depleted nitrogen-containing ring” indicates a C_1 - C_{60} heterocyclic group having at least one $*-N=*$ moiety as a ring-forming moiety.

[0256] For example, the “ π electron-depleted nitrogen-containing ring” may be i) a 5-membered to 7-membered heteromonocyclic group having at least one $*-N=*$ moiety, ii) a heteropolycyclic group in which two or more 5-membered to 7-membered heteromonocyclic groups each having at least one $*-N=*$ moiety are condensed with each other, or iii) a heteropolycyclic group in which at least one of 5-membered to 7-membered heteromonocyclic groups, each having at least one $*-N=*$ moiety, is condensed with at least one C_5 - C_{60} carbocyclic group.

[0257] Examples of the π electron-depleted nitrogen-containing ring include an imidazole, a pyrazole, a thiazole, an isothiazole, an oxazole, an isoxazole, a pyridine, a pyrazine, a pyrimidine, a pyridazine, an indazole, a purine, a quinoline, an isoquinoline, a benzoquinoline, a phthalazine, a naphthyridine, a quinoxaline, a quinazoline, a cinnoline, a phenanthridine, an acridine, a phenanthroline, a phenazine, a benzimidazole, an isobenzothiazole, a benzoxazole, an isobenzoxazole, a triazole, a tetrazole, an oxadiazole, a triazine, thiadiazol, an imidazopyridine, an imidazopyrimidine, and an azacarbazole, but are not limited thereto.

[0258] For example, the electron transport region may include a compound represented by Formula 601:



[0259] In Formula 601,

[0260] Ar_{601} may be a substituted or unsubstituted C_5 - C_{60} carbocyclic group or a substituted or unsubstituted C_1 - C_{60} heterocyclic group,

[0261] $xe11$ may be 1, 2, or 3,

[0262] L_{601} may be 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,

[0263] $xe1$ may be an integer selected from 0 to 5,

[0264] R_{601} 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, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, $-\text{Si}(Q_{601})(Q_{602})(Q_{603})$, $-\text{C}(=\text{O})(Q_{601})$, $-\text{S}(=\text{O})_2(Q_{601})$, and $-\text{P}(=\text{O})(Q_{601})(Q_{602})$,

[0265] Q_{601} to Q_{603} may each independently be a C_1 - C_{10} alkyl group, a C_1 - C_{10} alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, or a naphthyl group, and

[0266] $xe21$ may be an integer from 1 to 5.

[0267] In one or more embodiments, at least one of $Ar_{601}(s)$ in the number of $xe11$ and/or at least one of $R_{601}(s)$ in the number of $xe21$ may include the π electron-depleted nitrogen-containing ring.

[0268] In one or more embodiments, ring Ar_{601} in Formula 601 may be selected from:

[0269] a benzene group, a naphthalene group, a fluorene group, a spiro-bifluorene 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, an indenoanthracene group, a dibenzofuran group, a dibenzothiophene group, a carbazole group, an imidazole group, a pyrazole group, a thiazole group, an isothiazole group, an oxazole group, an isoxazole group, a pyridine group, a pyrazine group, a pyrimidine group, a pyridazine group, an indazole group, a purine group, a quinoline group, an isoquinoline group, a benzoquinoline group, a phthalazine group, a naphthyridine group, a quinoxaline group, a quinazoline group, a cinnoline group, a phenanthridine group, an acridine group, a phenanthroline group, a phenazine group, a benzimidazole group, an isobenzothiazole group, a benzoxazole group, an isobenzoxazole group, a triazole group, a tetrazole group, an oxadiazole group, a triazine group, thiadiazol group, an imidazopyridine group, an imidazopyrimidine group, and an azacarbazole group; and

[0270] a benzene group, a naphthalene group, a fluorene group, a spiro-bifluorene 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, an indenoanthracene group, a dibenzofuran group, a dibenzothiophene group, a carbazole group, an imidazole group, a pyrazole group, a thiazole group, an isothiazole group, an oxazole group, an isoxazole group, a pyridine group, a pyrazine group, a pyrimidine group, a pyridazine group, an indazole group, a purine group, a quinoline group, an isoquinoline group, a benzoquinoline group, a phthalazine group, a naphthyridine group, a quinoxaline group, a quinazoline group, a cinnoline group, a phenanthridine group, an acridine group, a phenanthroline group, a phenazine group, a benzimidazole group, an isobenzothiazole group, a benzoxazole group, an isobenzoxazole group, a triazole group, a tetrazole group, an oxadiazole group, a triazine group, thiadiazol group, an imidazopyridine group, an imidazopyrimidine group, and an azacarbazole group, each substituted with at

least one selected from deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, —Si(Q₃₁)(Q₃₂)(Q₃₃), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂), and

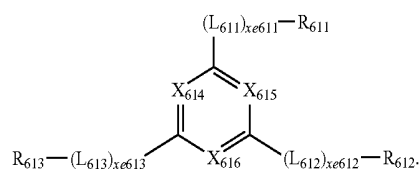
[0271] Q₃₁ to Q₃₃ may each independently be selected from a C₁-C₁₀ alkyl group, a C₁-C₁₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group.

[0272] When xe11 in Formula 601 is two or more, two or more Ar₆₀₁(s) may be linked via a single bond.

[0273] In one or more embodiments, Ar₆₀₁ in Formula 601 may be an anthracene group.

[0274] In one or more embodiments, a compound represented by Formula 601 may be represented by Formula 601-1:

<Formula 601-1>



[0275] In Formula 601-1,

[0276] X₆₁₄ may be N or C(R₆₁₄), X₆₁₅ may be N or C(R₆₁₅), X₆₁₆ may be N or C(R₆₁₆), and at least one selected from X₆₁₄ to X₆₁₆ may be N,

[0277] L₆₁₁ to L₆₁₃ may each independently be the same as described in connection with L₆₀₁,

[0278] xe611 to xe613 may each independently be the same as described in connection with xe1,

[0279] R₆₁₁ to R₆₁₃ may each independently be the same as described in connection with R₆₀₁, and

[0280] R₆₁₄ to R₆₁₆ may each independently be selected from hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group.

[0281] In one or more embodiments, L₆₀₁ and L₆₁₁ to L₆₁₃ in Formulae 601 and 601-1 may each independently be selected from:

[0282] a phenylene group, a naphthylene group, a fluorenylene group, a spiro-bifluorenylene group, a benzofluorenylene group, a dibenzofluorenylene group, a phenanthrenylene group, an anthracenylylene group, a fluoranthrenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a perylenylene group, a pentaphenylylene group, a hexacenylylene group, a pentacenylylene group, a thiophenylylene group, a furanylylene group, a carbazolylylene group, an indolylylene group, an isoindolylylene group, a benzofuranylylene group, a benzothiophenylylene group, a dibenzofuranylylene group, a dibenzothiophenylylene group, a benzocarbazolylylene group, a dibenzocarbazolylylene

group, a dibenzosilolylylene group, a pyridinylylene group, an imidazolylylene group, a pyrazolylylene group, a thiazolylylene group, an isothiazolylylene group, an oxazolylylene group, an isoxazolylylene group, a thiadiazolylylene group, an oxadiazolylylene group, a pyrazinylylene group, a pyrimidinylylene group, a pyridazinylylene group, a triazinylylene group, a quinolinylene group, an isoquinolinylene group, a benzoquinolinylene group, a phthalazinylylene group, a naphthyridinylylene group, a quinoxalinylylene group, a quinazolinylylene group, a cinolinylene group, a phenanthridinylylene group, an acridinylylene group, a phenanthrolinylylene group, a phenazinylylene group, a benzimidazolylylene group, an isobenzothiazolylylene group, a benzoxazolylylene group, an isobenzoxazolylylene group, a triazolylylene group, a tetrazolylylene group, an imidazopyridinylylene group, an imidazopyrimidinylylene group, and an azacarbazolylylene group; and

[0283] a phenylene group, a naphthylene group, a fluorenylylene group, a spiro-bifluorenylylene group, a benzofluorenylylene group, a dibenzofluorenylylene group, a phenanthrenylene group, an anthracenylylene group, a fluoranthrenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a perylenylene group, a pentaphenylylene group, a hexacenylylene group, a pentacenylylene group, a thiophenylylene group, a furanylylene group, a carbazolylylene group, an indolylylene group, an isoindolylylene group, a benzofuranylylene group, a benzothiophenylylene group, a dibenzofuranylylene group, a dibenzothiophenylylene group, a benzocarbazolylylene group, a dibenzocarbazolylylene group, a dibenzosilolylylene group, a pyridinylylene group, an imidazolylylene group, a pyrazolylylene group, a thiazolylylene group, an isothiazolylylene group, an oxazolylylene group, an isoxazolylylene group, a thiadiazolylylene group, an oxadiazolylylene group, a pyrazinylylene group, a pyrimidinylylene group, a pyridazinylylene group, a triazinylylene group, a quinolinylene group, an isoquinolinylene group, a benzoquinolinylene group, a phthalazinylylene group, a naphthyridinylylene group, a quinoxalinylylene group, a quinazolinylylene group, a cinolinylene group, a phenanthridinylylene group, an acridinylylene group, a phenanthrolinylylene group, a phenazinylylene group, a benzimidazolylylene group, an isobenzothiazolylylene group, a benzoxazolylylene group, an isobenzoxazolylylene group, a triazolylylene group, a tetrazolylylene group, an imidazopyridinylylene group, an imidazopyrimidinylylene group, and an azacarbazolylylene group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthrenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl

group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, a pyridinyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a thiadiazolyl group, an oxadiazolyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a triazinyl 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 phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzimidazolyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, a tetrazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and an azacarbazolyl group, but are not limited thereto.

[0284] but embodiments of the present disclosure are not limited thereto.

[0285] In one or more embodiments, xe1 and xe11 to xe613 in Formulae 601 and 601-1 may each independently be 0, 1, or 2.

[0286] In one or more embodiments, R₆₀₁ and R₆₁₁ to R₆₁₃ in Formula 601 and 601-1 may each independently be selected from:

[0287] a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, a pyridinyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a thiadiazolyl group, an oxadiazolyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a triazinyl 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 phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzimidazolyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, a tetrazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and an azacarbazolyl group;

[0288] a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a

pyrenyl group, a chrysenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, a pyridinyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a thiadiazolyl group, an oxadiazolyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a triazinyl 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 phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzimidazolyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, a tetrazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and an azacarbazolyl group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a thiophenyl group, a furanyl group, a carbazolyl group, an indolyl group, an isoindolyl group, a benzofuranyl group, a benzothiophenyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a dibenzosilolyl group, a pyridinyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a thiadiazolyl group, an oxadiazolyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a triazinyl 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 phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzimidazolyl group, an isobenzothiazolyl group, a benzoxazolyl group, an isobenzoxazolyl group, a triazolyl group, a tetrazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and an azacarbazolyl group; and

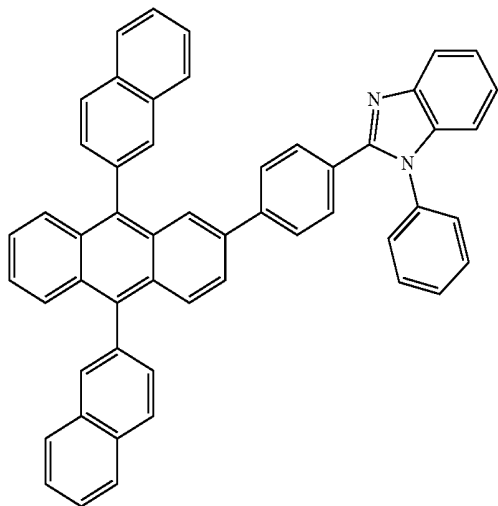
[0289] —S(=O)₂(Q₆₀₁), and —P(=O)(Q₆₀₁)(Q₆₀₂); and

[0290] Q₆₀₁ and Q₆₀₂ may be the same as described above.

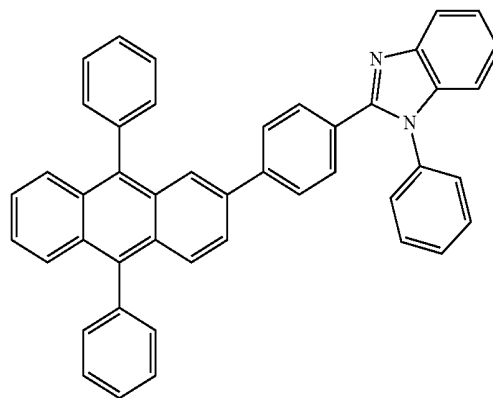
[0291] The electron transport region may include at least one compound selected from Compounds ET1 to ET36, but embodiments of the present disclosure are not limited thereto:

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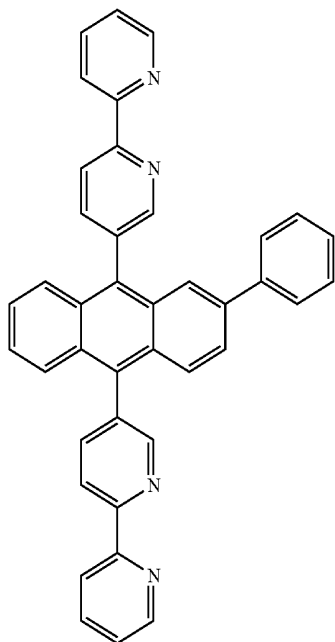
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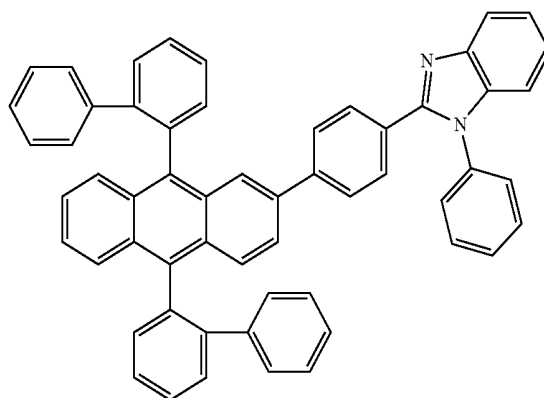
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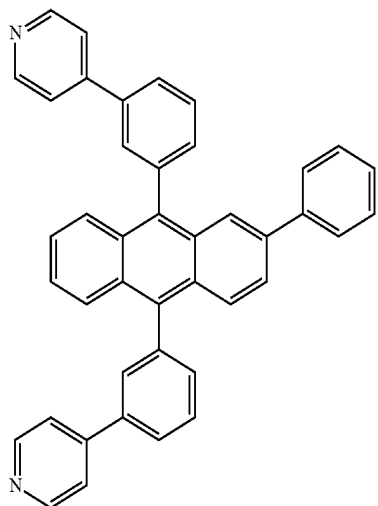
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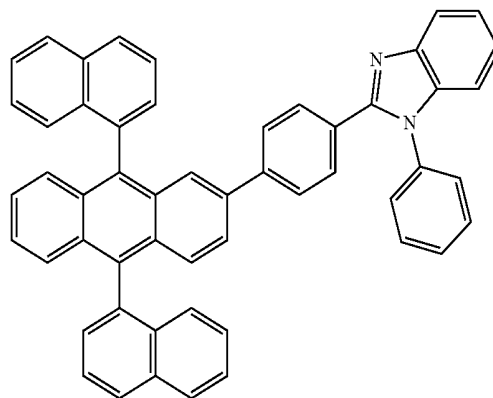
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ET3

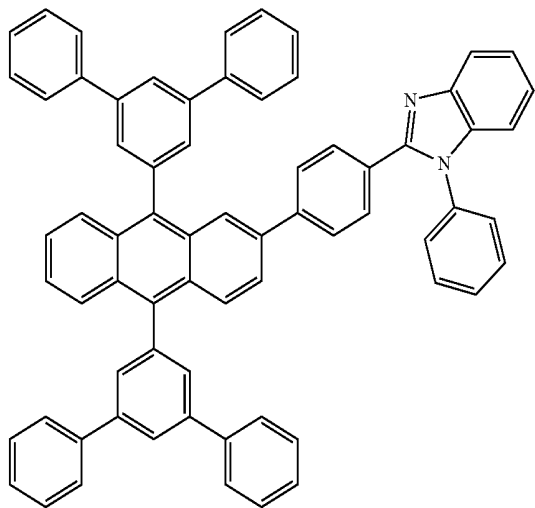


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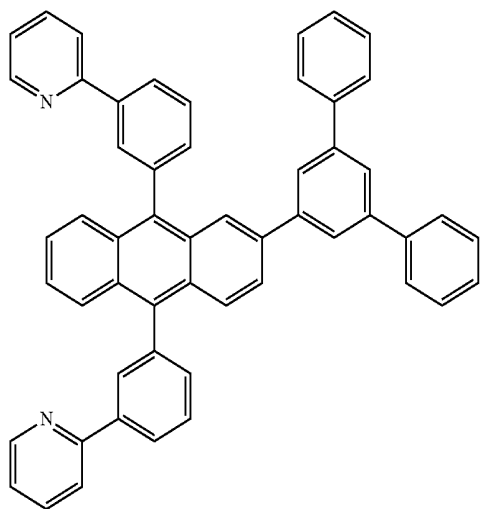


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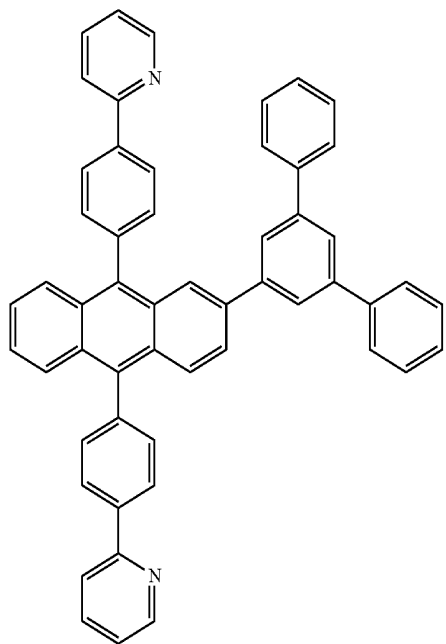
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ET8

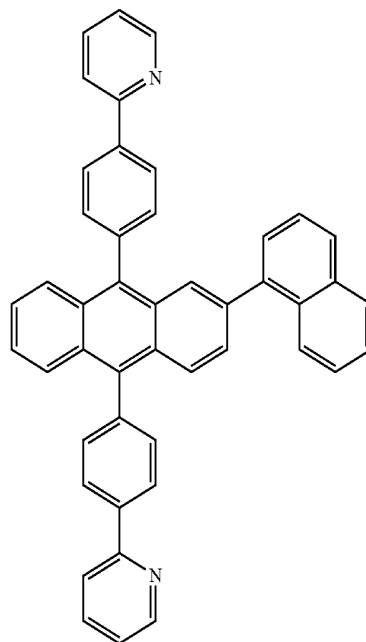


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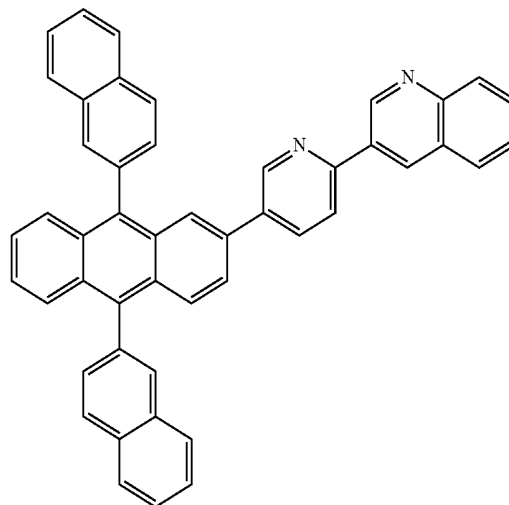


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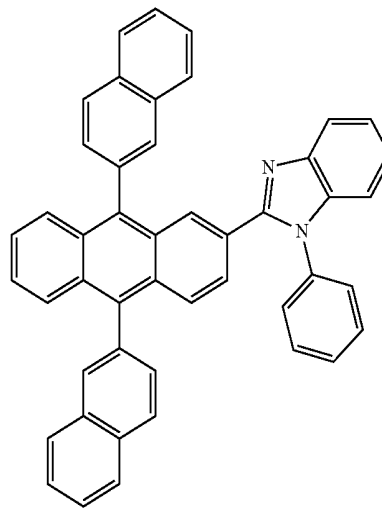
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ET11



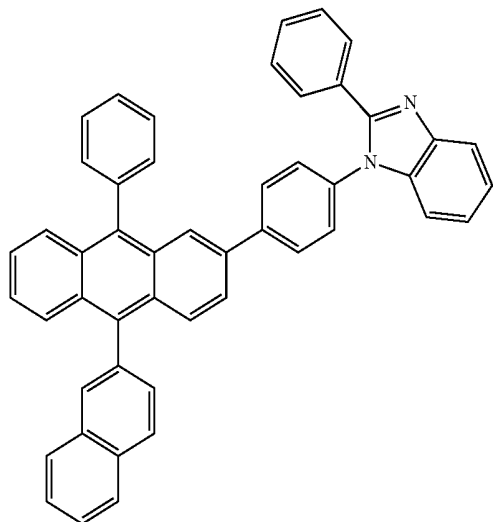
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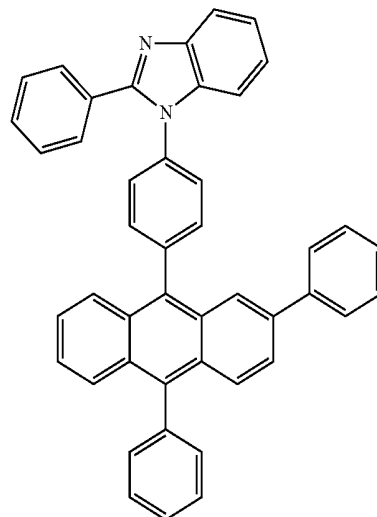
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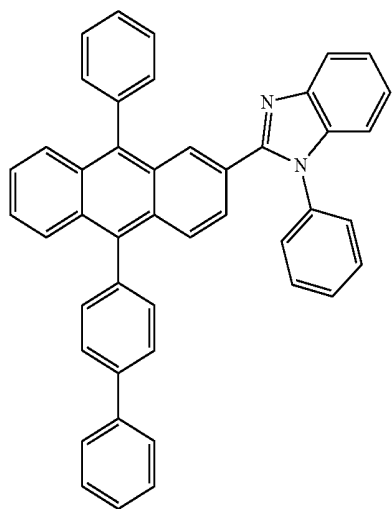
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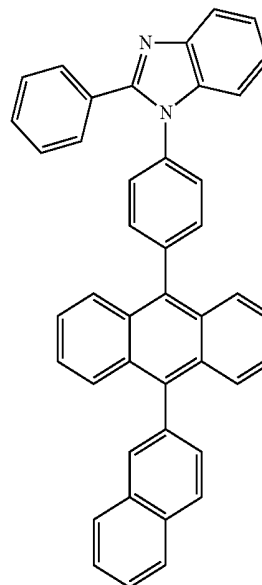
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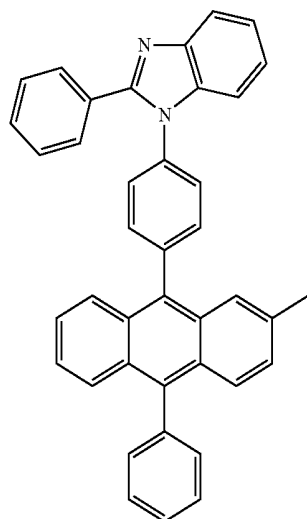
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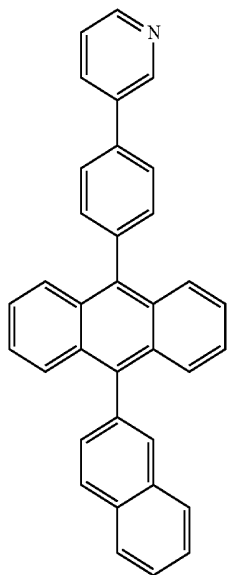
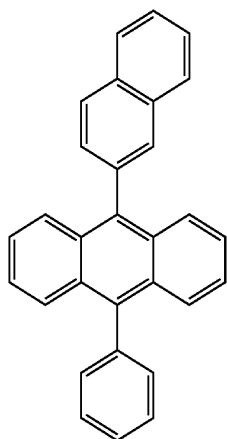
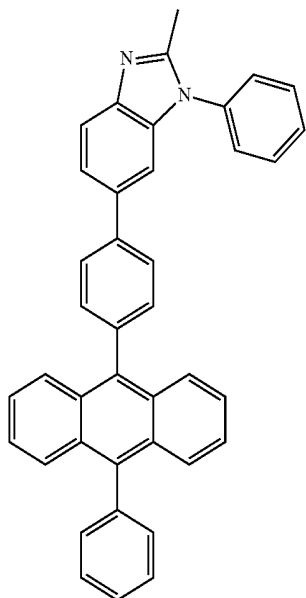
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ET15

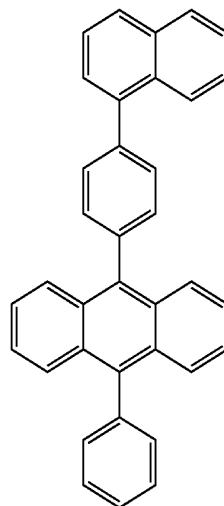


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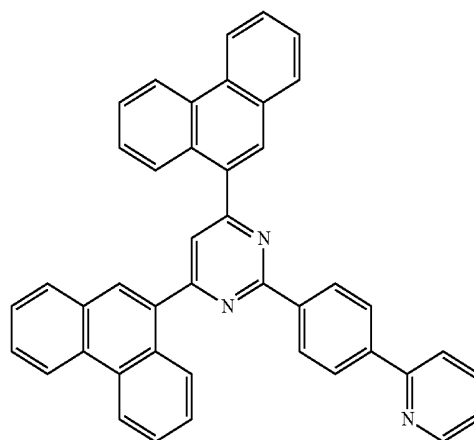
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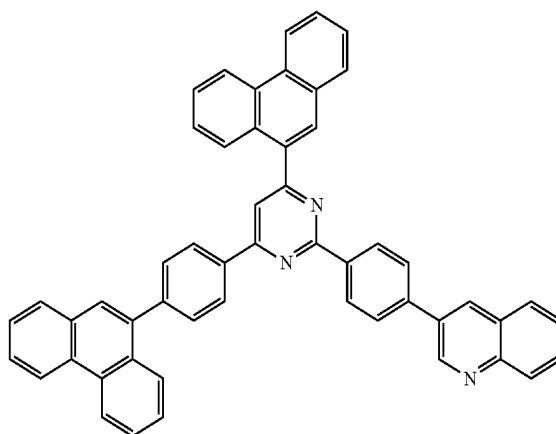
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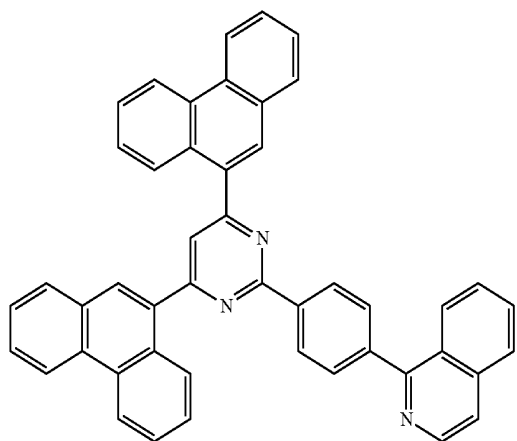
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ET23

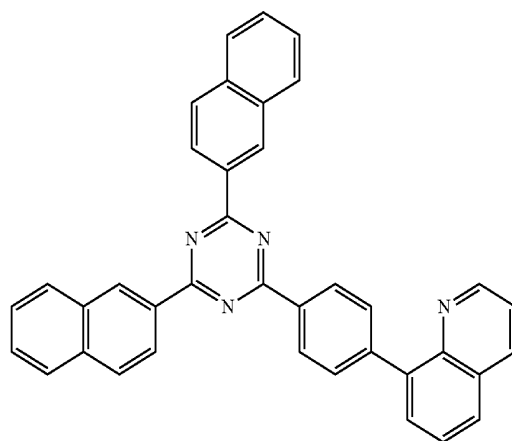
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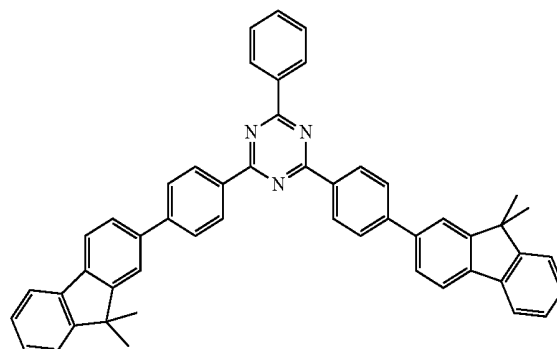
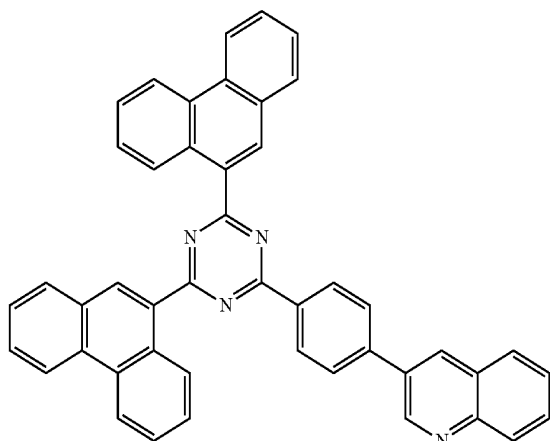
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ET27



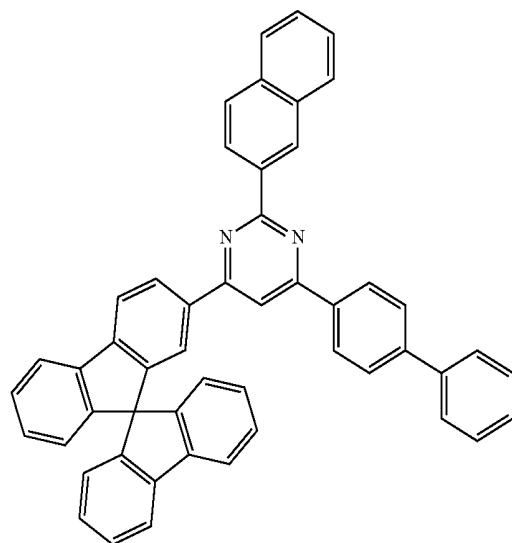
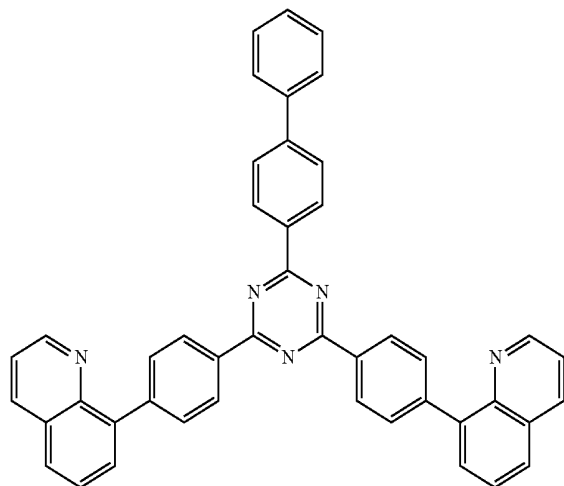
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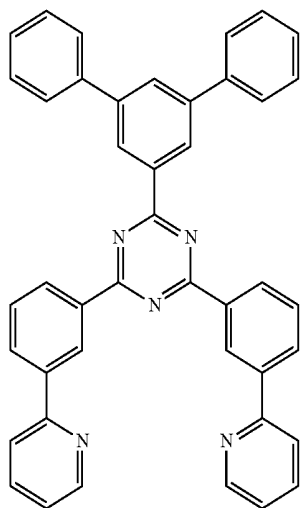


ET26

ET29

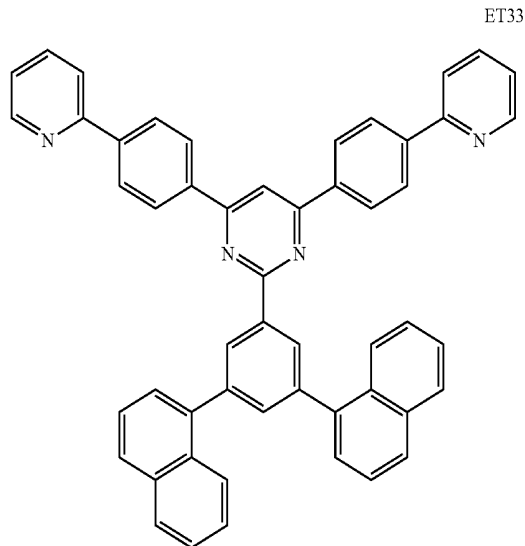


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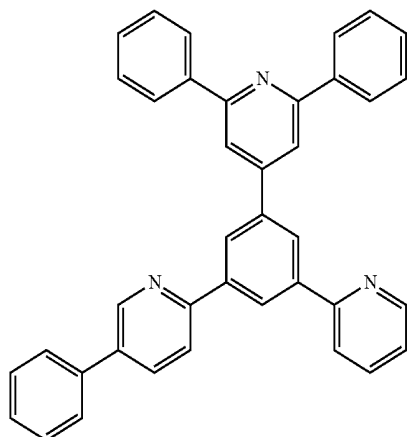


ET30

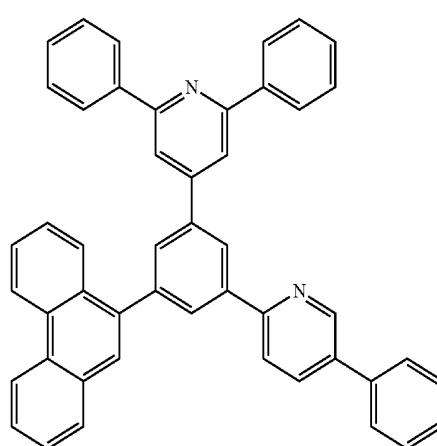
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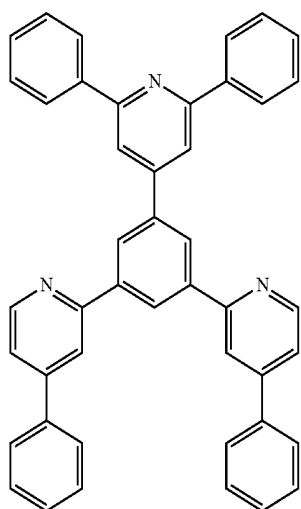
ET33



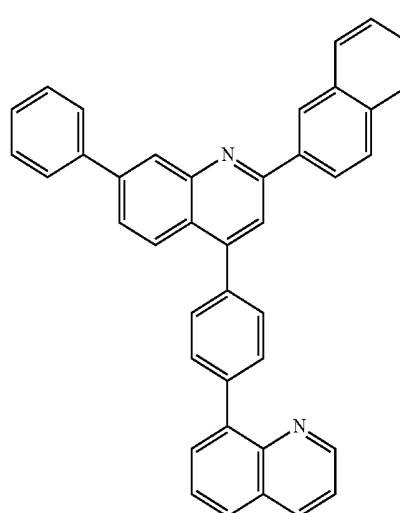
ET31



ET34

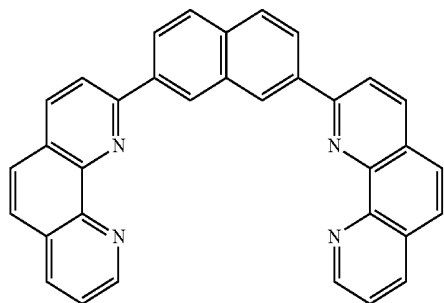


ET32



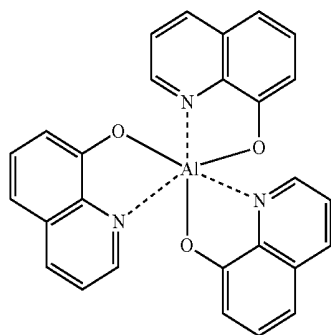
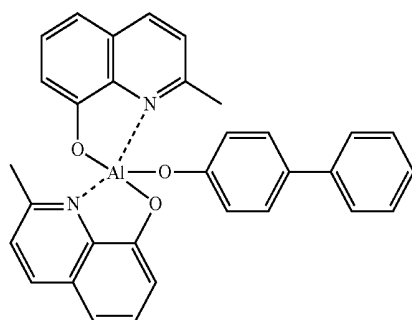
ET35

-continued

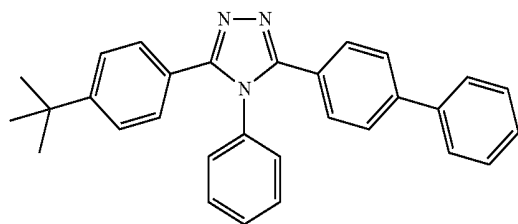


ET36

[0292] In one or more embodiments, the electron transport region may include at least one selected from 2,9-dimethyl-4,7-diphenyl-1,10-phenanthroline (BCP), 4,7-diphenyl-1,10-phenanthroline (Bphen), Alq₃, BAlq, 3-(biphenyl-4-yl)-5-(4-tert-butylphenyl)-4-phenyl-4H-1,2,4-triazole (TAZ), and NTAZ.

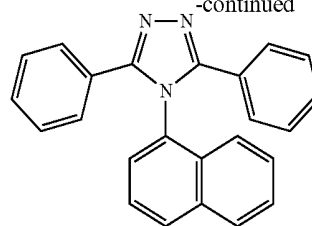
Alq₃

BAlq



TAZ

-continued



NTAZ

[0293] In one or more embodiments, the electron transport region may include a phosphine oxide-containing compound (for example, TSPO1 used in the following examples or the like), but embodiments of the present disclosure are not limited thereto. In one or more embodiments, the phosphine oxide-containing compound may be used in a hole blocking layer in the electron transport region, but embodiments of the present disclosure are not limited thereto.

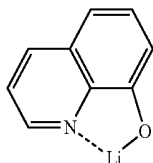
[0294] The thickness of the buffer layer, the hole blocking layer, or the electron controlling layer may each independently be in a range of about 20 Å to about 1,000 Å, for example, about 30 Å to about 300 Å. When the thicknesses of the buffer layer, the hole blocking layer, and the electron control layer are within these ranges, the electron blocking layer may have excellent electron blocking characteristics or electron control characteristics without a substantial increase in driving voltage.

[0295] A thickness of the electron transport layer may be in a range of about 100 Å to about 1,000 Å, for example, about 150 Å to about 500 Å. When the thickness of the electron transport layer is within the range described above, the electron transport layer may have satisfactory electron transport characteristics without a substantial increase in driving voltage.

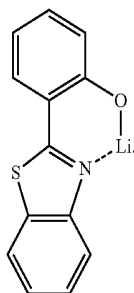
[0296] The electron transport region (for example, the electron transport layer in the electron transport region) may further include, in addition to the materials described above, a metal-containing material.

[0297] The metal-containing material may include at least one selected from alkali metal complex and alkaline earth-metal complex. The alkali metal complex may include a metal ion selected from a Li ion, a Na ion, a K ion, a Rb ion, and a Cs ion, and the alkaline earth-metal complex may include a metal ion selected from a Be ion, a Mg ion, a Ca ion, a Sr ion, and a Ba ion. A ligand coordinated with the metal ion of the alkali metal complex or the alkaline earth-metal complex may be selected from a hydroxy quinoline, a hydroxy isoquinoline, a hydroxy benzoquinoline, a hydroxy acridine, a hydroxy phenanthridine, a hydroxy phenyl oxazole, a hydroxy phenylthiazole, a hydroxy diphenyl oxadiazole, a hydroxy diphenylthiadiazol, a hydroxy phenylpyridine, a hydroxy phenylbenzimidazole, a hydroxy phenylbenzothiazole, a bipyridine, a phenanthroline, and a cyclopentadiene, but embodiments of the present disclosure are not limited thereto.

[0298] For example, 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

[0299] The electron transport region may include an electron injection layer that facilitates injection of electrons from the second electrode **190**. The electron injection layer may directly contact the second electrode **190**.

[0300] The electron injection layer may have i) a single-layered structure including a single layer including a single material, ii) a single-layered structure including a single layer including a plurality of different materials, or iii) a multi-layered structure having a plurality of layers including a plurality of different materials.

[0301] The electron injection layer may include an alkali metal, an alkaline earth metal, a rare earth metal, an alkali metal compound, an alkaline earth-metal compound, a rare earth metal compound, an alkali metal complex, an alkaline earth-metal complex, a rare earth metal complex, or any combinations thereof.

[0302] The alkali metal may be selected from Li, a Na, K, Rb, and Cs. In one or more embodiments, the alkali metal may be Li, a Na, or Cs. In one or more embodiments, the alkali metal may be Li or Cs, but embodiments of the present disclosure are not limited thereto.

[0303] The alkaline earth metal may be selected from Mg, Ca, Sr, and Ba.

[0304] The rare earth metal may be selected from Sc, Y, Ce, Tb, Yb, and Gd.

[0305] The alkali metal compound, the alkaline earth-metal compound, and the rare earth metal compound may be selected from oxides and halides (for example, fluorides, chlorides, bromides, or iodides) of the alkali metal, the alkaline earth-metal, and the rare earth metal.

[0306] The alkali metal compound may be selected from alkali metal oxides, such as Li_2O , Cs_2O , or K_2O , and alkali metal halides, such as LiF, a NaF, CsF, KF, LiI, a NaI, CsI, or KI. In one or more embodiments, the alkali metal compound may be selected from LiF, Li_2O , a NaF, LiI, a NaI, CsI, and KI, but embodiments of the present disclosure are not limited thereto.

[0307] The alkaline earth-metal compound may be selected from alkaline earth-metal oxides, such as BaO, SrO, CaO, $\text{Ba}_x\text{Sr}_{1-x}\text{O}$ ($0 < x < 1$), or $\text{Ba}_x\text{Ca}_{1-x}\text{O}$ ($0 < x < 1$). In one or more embodiments, the alkaline earth-metal compound may be selected from BaO, SrO, and CaO, but embodiments of the present disclosure are not limited thereto.

[0308] The rare earth metal compound may be selected from YbF_3 , ScF_3 , ScO_3 , Y_2O_3 , Ce_2O_3 , GdF_3 , and TbF_3 . In one or more embodiments, the rare earth metal compound may be selected from YbF_3 , ScF_3 , TbF_3 , YbI_3 , ScI_3 , and TbI_3 , but embodiments of the present disclosure are not limited thereto.

[0309] The alkali metal complex, the alkaline earth-metal complex, and the rare earth metal complex may include an ion of alkali metal, alkaline earth-metal, and rare earth metal as described above, and a ligand coordinated with a metal ion of the alkali metal complex, the alkaline earth-metal complex, or the rare earth metal complex may be selected from hydroxy quinoline, hydroxy isoquinoline, hydroxy benzoquinoline, hydroxy acridine, hydroxy phenanthridine, hydroxy phenyl oxazole, hydroxy phenylthiazole, hydroxy diphenyl oxadiazole, hydroxy diphenylthiadiazol, hydroxy phenylpyridine, hydroxy phenylbenzimidazole, hydroxy phenylbenzothiazole, bipyridine, phenanthroline, and cyclopentadiene, but embodiments of the present disclosure are not limited thereto.

[0310] The electron injection layer may consist of an alkali metal, an alkaline earth metal, a rare earth metal, an alkali metal compound, an alkaline earth-metal compound, a rare earth metal compound, an alkali metal complex, an alkaline earth-metal complex, a rare earth metal complex, or any combinations thereof, as described above. In one or more embodiments, the electron injection layer may further include an organic material.

[0311] When the electron injection layer further includes an organic material, an alkali metal, an alkaline earth metal, a rare earth metal, an alkali metal compound, an alkaline earth-metal compound, a rare earth metal compound, an alkali metal complex, an alkaline earth-metal complex, a rare earth metal complex, or any combinations thereof may be homogeneously or non-homogeneously dispersed in a matrix including the organic material.

[0312] A thickness of the electron injection layer may be in a range of about 1 Å to about 100 Å, for example, about 3 Å to about 90 Å. When the thickness of the electron injection layer is within the range described above, the electron injection layer may have satisfactory electron injection characteristics without a substantial increase in driving voltage.

[Second Electrode **190**]

[0313] 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 combination thereof.

[0314] The second electrode **190** may include at least one selected from lithium (Li), silver (Si), magnesium (Mg), aluminum (Al), aluminum-lithium (Al—Li), calcium (Ca), magnesium-indium (Mg—In), magnesium-silver (Mg—Ag), ITO, and IZO, but embodiments of the present disclosure are not limited thereto. The second electrode **190** may be a transmissive electrode, a semi-transmissive electrode, or a reflective electrode.

[0315] The second electrode **190** may have a single-layered structure, or a multi-layered structure including two or more layers.

[0316] Layers constituting the hole transport region, an emission layer, and layers constituting the electron transport region may be formed in a certain region by using one or more suitable methods selected from vacuum deposition, spin coating, casting, Langmuir-Blodgett (LB) deposition, ink-jet printing, laser-printing, and laser-induced thermal imaging.

[0317] When the respective layers of the hole transport region, the emission layer, and the respective layers of the electron transport region are formed by deposition, the 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 by taking into account a material for forming a layer to be deposited, and the structure of a layer to be formed.

[0318] When layers constituting the hole transport region, an emission layer, and layers constituting the electron transport region are formed by spin coating, the spin coating may be performed at a coating speed of about 2,000 rpm to about 5,000 rpm and at a heat treatment temperature of about 80° C. to 200° C., depending on a material to be included in a layer and the structure of each layer to be formed.

[General Definition of Substituents]

[0319] The term “C₁-C₆₀ alkyl group” as used herein refers to a linear or branched saturated aliphatic hydrocarbon monovalent group having 1 to 60 carbon atoms, and non-limiting examples thereof include 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. The term “C₁-C₆₀ alkylene group” as used herein refers to a divalent group having the same structure as the C₁-C₆₀ alkyl group.

[0320] The term “C₂-C₆₀ alkenyl group” as used herein refers to a hydrocarbon group formed by substituting at least one carbon-carbon double bond in the middle or at the terminus of the C₂-C₆₀ alkyl group, and non-limiting examples thereof include an ethenyl group, a propenyl group, and a butenyl group. The term “C₂-C₆₀ alkenylene group” as used herein refers to a divalent group having the same structure as the C₂-C₆₀ alkenyl group.

[0321] The term “C₂-C₆₀ alkynyl group” as used herein refers to a hydrocarbon group formed by substituting at least one carbon-carbon triple bond in the middle or at the terminus of the C₂-C₆₀ alkyl group, and non-limiting examples thereof include an ethynyl group, and a propynyl group. The term “C₂-C₆₀ alkynylene group” as used herein refers to a divalent group having the same structure as the C₂-C₆₀ alkynyl group.

[0322] The term “C₁-C₆₀ alkoxy group” as used herein refers to a monovalent group represented by —OA₁₀₁ (wherein A₁₀₁ is the C₁-C₆₀ alkyl group), and non-limiting examples thereof include a methoxy group, an ethoxy group, and an isopropoxy group.

[0323] The term “C₃-C₁₀ cycloalkyl group” as used herein refers to a monovalent saturated hydrocarbon monocyclic group having 3 to 10 carbon atoms, and non-limiting examples thereof include a cyclopropyl group, a cyclobutyl group, a cyclopentyl group, a cyclohexyl group, and a cycloheptyl group. The term “C₃-C₁₀ cycloalkylene group” as used herein refers to a divalent group having the same structure as the C₃-C₁₀ cycloalkyl group.

[0324] The term “C₁₀-C₁₀ heterocycloalkyl group” as used herein refers to a monovalent saturated monocyclic group having at least one heteroatom selected from N, O, Si, P, and S as a ring-forming atom and 1 to 10 carbon atoms, and non-limiting examples thereof include a 1,2,3,4-oxatriazolidynyl group, a tetrahydrofuranyl group, and a tetrahydrothiophenyl group. The term “C₁-C₁₀ heterocycloalkylene group” as used herein refers to a divalent group having the same structure as the C₁-C₁₀ heterocycloalkyl group.

[0325] The term “C₃-C₁₀ cycloalkenyl group” as used herein refers to a monovalent monocyclic group that has 3 to 10 carbon atoms and at least one carbon-carbon double bond in the ring thereof and no aromaticity, and non-limiting examples thereof include a cyclopentenyl group, a cyclohexenyl group, and a cycloheptenyl group. The term “C₃-C₁₀ cycloalkenylene group” as used herein refers to a divalent group having the same structure as the C₃-C₁₀ cycloalkenyl group.

[0326] The term “C₁-C₁₀ heterocycloalkenyl group” as used herein refers to a monovalent monocyclic group that has at least one heteroatom selected from N, O, Si, P, and S as a ring-forming atom, 1 to 10 carbon atoms, and at least one carbon-carbon double bond in its ring. Non-limiting examples of the C₁-C₁₀ heterocycloalkenyl group include a 4,5-dihydro-1,2,3,4-oxatriazolyl group, a 2,3-dihydrofuranyl group, and a 2,3-dihydrothiophenyl group. The term “C₁-C₁₀ heterocycloalkenylene group” as used herein refers to a divalent group having the same structure as the C₁-C₁₀ heterocycloalkenyl group.

[0327] The term “C₆-C₆₀ aryl group” as used herein refers to a monovalent group having a carbocyclic aromatic system having 6 to 60 carbon atoms, and the term “C₆-C₆₀ arylene group” as used herein refers to a divalent group having a carbocyclic aromatic system having 6 to 60 carbon atoms. Non-limiting examples of the C₆-C₆₀ aryl group include 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.

[0328] The term “C₁-C₆₀ heteroaryl group” as used herein refers to a monovalent group having a heterocyclic aromatic system that has at least one heteroatom selected from N, O, Si, P, and S as a ring-forming atom, in addition to 1 to 60 carbon atoms. The term “C₁-C₆₀ heteroarylene group” as used herein refers to a divalent group having a heterocyclic aromatic system that has at least one heteroatom selected from N, O, Si, P, and S as a ring-forming atom, in addition to 1 to 60 carbon atoms. Non-limiting examples of the C₁-C₆₀ heteroaryl group include 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.

[0329] The term “C₆-C₆₀ aryloxy group” as used herein refers to —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).

[0330] The term “monovalent non-aromatic condensed polycyclic group” as used herein refers to a monovalent group (for example, having 8 to 60 carbon atoms) having two or more rings condensed with each other, only carbon atoms as ring-forming atoms, and no aromaticity in its entire

molecular structure. A detailed example of the monovalent non-aromatic condensed polycyclic group is a fluorenyl group. The term “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.

[0331] The term “monovalent non-aromatic condensed heteropolycyclic group” as used herein refers to a monovalent group (for example, having 1 to 60 carbon atoms) having two or more rings condensed to each other, at least one heteroatom selected from N, O, Si, P, and S, other than carbon atoms, as a ring-forming atom, and no aromaticity in its entire molecular structure. An example of the monovalent non-aromatic condensed heteropolycyclic group is a carbazolyl group. The term “divalent non-aromatic condensed heteropolycyclic group” as used herein refers to a divalent group having the same structure as the monovalent non-aromatic condensed heteropolycyclic group.

[0332] The term “C₄-C₆₀ carbocyclic group” as used herein refers to a monocyclic or polycyclic group having 4 to 60 carbon atoms in which a ring-forming atom is a carbon atom only. The C₄-C₆₀ carbocyclic group may be an aromatic carbocyclic group or a non-aromatic carbocyclic group. The C₄-C₆₀ carbocyclic group may be a ring, such as benzene, a monovalent group, such as a phenyl group, or a divalent group, such as a phenylene group. In one or more embodiments, depending on the number of substituents connected to the C₄-C₆₀ carbocyclic group, the C₄-C₆₀ carbocyclic group may be a trivalent group or a quadrivalent group.

[0333] The term “C₂-C₆₀ heterocyclic group” as used herein refers to a group having the same structure as the C₄-C₆₀ carbocyclic group, except that as a ring-forming atom, at least one heteroatom selected from N, O, Si, P, and S is used in addition to carbon (the number of carbon atoms may be in a range of 1 to 60).

[0334] At least one substituent of the substituted C₄-C₆₀ carbocyclic group, substituted C₂-C₆₀ heterocyclic 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-C₁₀ 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:

[0335] deuterium (-D), -F, -Cl, -Br, -I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group;

[0336] 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 deuterium, -F, -Cl, -Br, -I, a hydroxyl group, a cyano group,

a nitro group, an amidino group, a hydrazino group, a hydrazono 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(Q₁₁)(Q₁₂)(Q₁₃), -N(Q₁₁)(Q₁₂), -B(Q₁₁)(Q₁₂), -C(=O)(Q₁₁), -S(=O)₂(Q₁₁), and -P(=O)(Q₁₁)(Q₁₂);

[0337] 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;

[0338] 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 deuterium, -F, -Cl, -Br, -I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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), -N(Q₂₁)(Q₂₂), -B(Q₂₁)(Q₂₂), -C(=O)(Q₂₁), -S(=O)₂(Q₂₁) and -P(=O)(Q₂₁)(Q₂₂); and

[0339] -Si(Q₃₁)(Q₃₂)(Q₃₃), -N(Q₃₁)(Q₃₂), -B(Q₃₁)(Q₃₂), -C(=O)(Q₃₁), -S(=O)₂(Q₃₁), and -P(=O)(Q₃₁)(Q₃₂); and

[0340] Q₁₁ to Q₁₃, Q₂₁ to Q₂₃ and Q₃₁ to Q₃₃ may each independently be selected from hydrogen, deuterium, -F, -Cl, -Br, -I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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, a monovalent non-aromatic condensed heteropolycyclic group, a C₁-C₆₀ alkyl group substituted with at least one selected from deuterium, -F, and a cyano group, a C₆-C₆₀ aryl group substituted with at least one selected from deuterium, -F, and a cyano group, a biphenyl group, and a terphenyl group.

[0341] The term “Ph”, as used herein, may refer to a phenyl group; the term “Me”, as used herein, may refer to a methyl group; the term “Et”, as used herein, may refer to an ethyl group; the terms “ter-Bu” or “But”, as used herein, may refer to a tert-butyl group; and the term “OMe” as used herein may refer to a methoxy group.

[0342] The term “biphenyl group” as used herein refers to “a phenyl group substituted with a phenyl group.” In other words, the “biphenyl group” is a substituted phenyl group having a C₆-C₆₀ aryl group as a substituent.

[0343] The term “a terphenyl group” as used herein refers to “a phenyl group substituted with a biphenyl group”. In other words, the “terphenyl group” is a phenyl group having, as a substituent, a C₆-C₆₀ aryl group substituted with a C₆-C₆₀ aryl group.

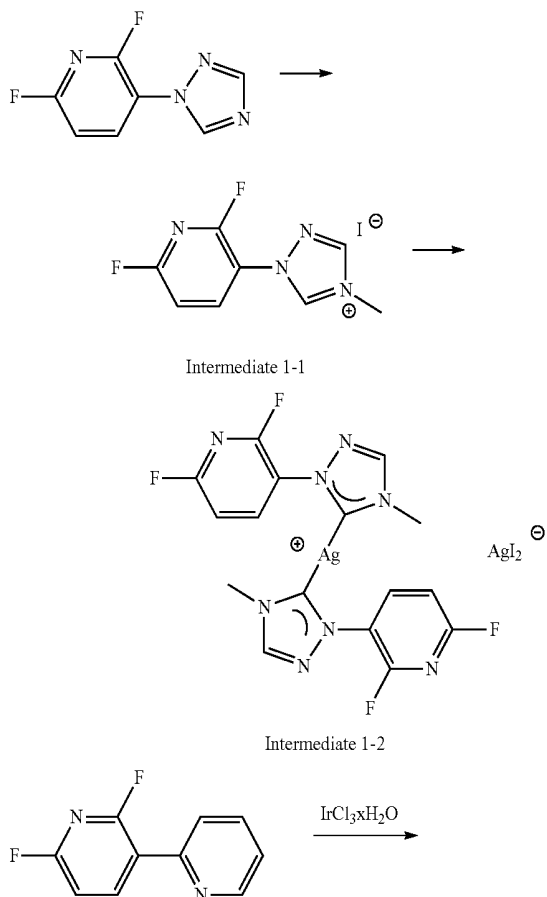
[0344] * and * used herein, unless defined otherwise, each refer to a binding site to a neighboring atom in a corresponding formula.

[0345] Hereinafter, a compound according to embodiments and an organic light-emitting device according to embodiments 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 refers to that an identical molar equivalent of B was used in place of A.

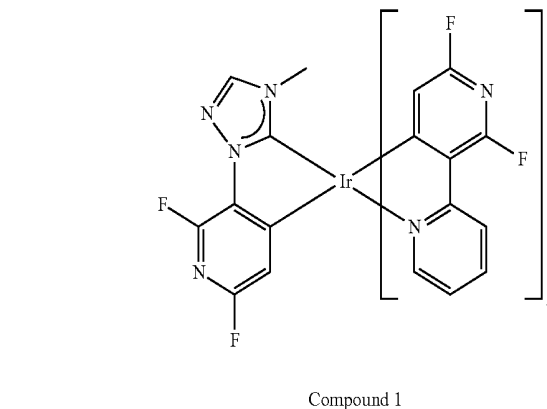
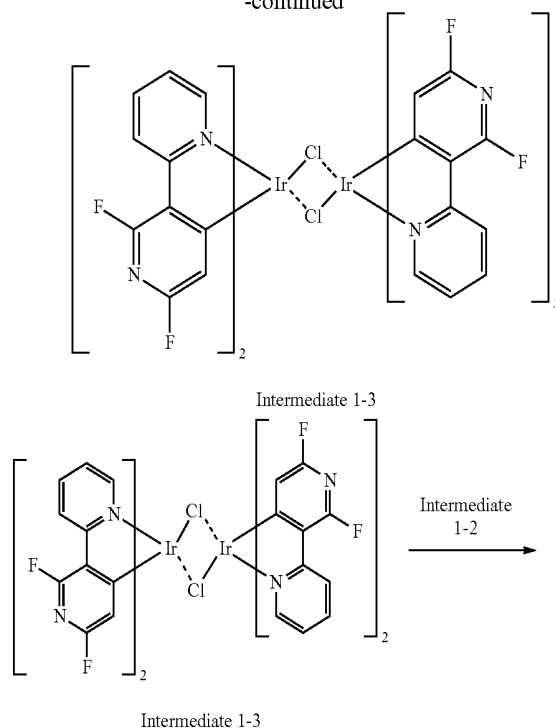
EXAMPLE

Synthesis Example 1: Synthesis of Compound 1

[0346]



-continued

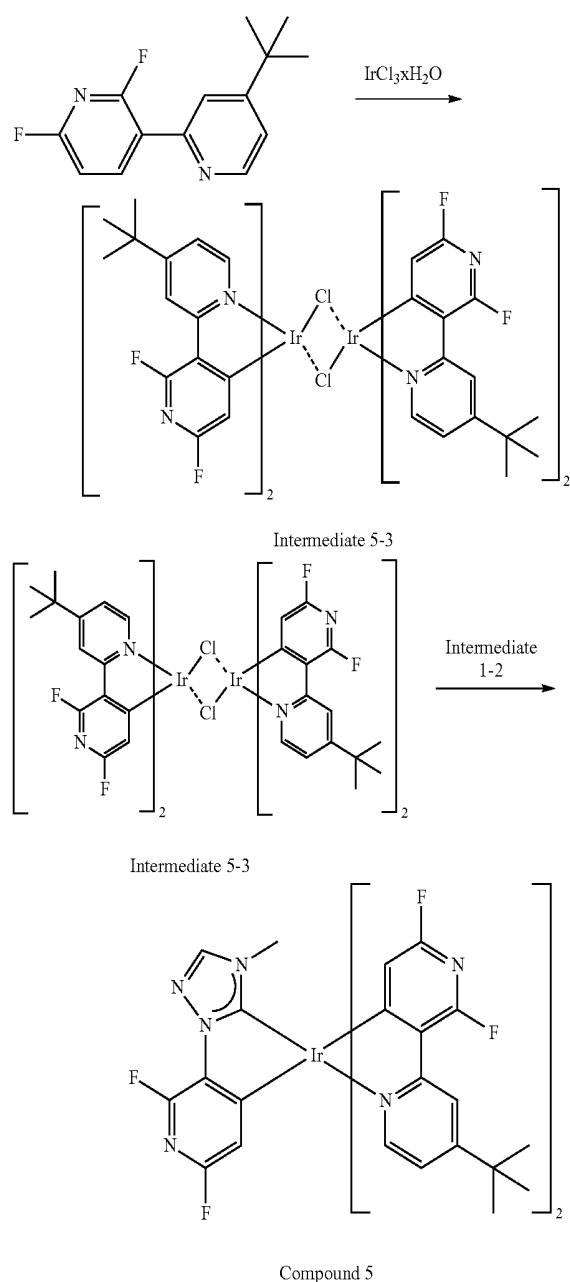


[0347] 2.00 g (10.98 mmol) of 2,6-difluoro-3-(1H-1,2,4-triazol-1-yl)pyridine and 3.42 mL (54.90 mmol) of iodomethane were mixed with 15 mL of tetrahydrofuran and then stirred at a temperature of 100° C. for 2 hours. The reaction mixture obtained therefrom was cooled to room temperature and a solvent was removed therefrom. The resultant was recrystallized by using diethyl ether and methanol to obtain Intermediate 1-1 (triazolium iodide salt). Intermediate 1-1 and 2.54 g (10.98 mmol) of silver (I) oxide were stirred in 15 mL of a dichloromethane solvent at room temperature for 6 hours. The solid obtained therefrom was filtered and a solvent was removed from a filtrate to obtain Intermediate 1-2 (silver carbene Intermediate). Then, Intermediate 1-2, 2',6'-difluoro-2,3'-bipyridine, and iridium (III) chloride hydrate react with one another in a 2-ethoxyethanol solvent at a temperature of 120° t for 6 hours to obtain 6.69 g (5.49 mmol) of Intermediate 1-3 (iridium dimer). Intermediate 1-3 was mixed with 30 mL of tetrahydrofuran and stirred at a temperature of 70° for 24 hours. The reaction mixture obtained therefrom was cooled to room temperature. An extraction process was performed thereon three

times by using dichloromethane and water. An organic layer obtained therefrom was dried by using magnesium sulfate and filtered by using celite, and column chromatography was performed thereon to synthesize 2.11 g (2.75 mmol, yield=25%) of Compound 1.

Synthesis Example 2: Synthesis of Compound 5

[0348]

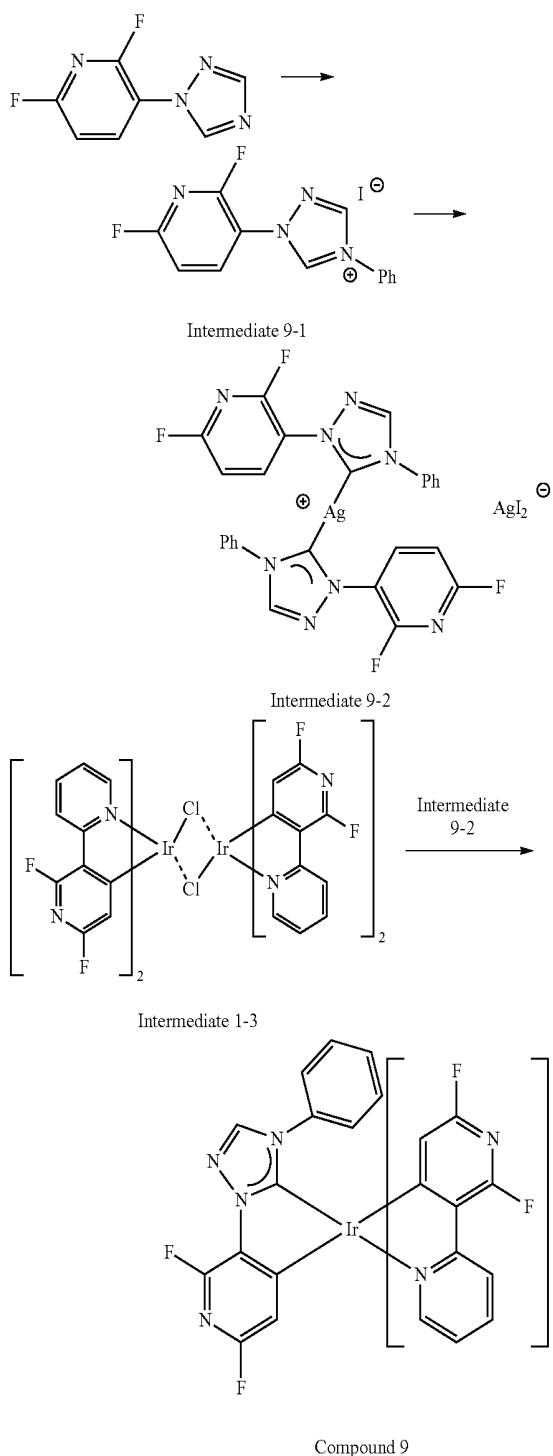


[0349] 7.93 g (5.49 mmol) of Intermediate 5-3 was obtained in the same manner as in the synthesis method of Intermediate 1-3 in Synthesis Example 1, except that 4-(tert-butyl)-2',6'-difluoro-2,3'-bipyridine was used instead of 2',6'-difluoro-2,3'-bipyridine in synthesizing Intermediate 1-3, and 1.75 g (1.98 mmol, yield=18%) of Compound 5 was synthesized in the same manner as in Synthesis of Com-

ound 1 in Synthesis Example 1, except that Intermediate 5-3 was used instead of Intermediate 1-3 in synthesizing Compound 1.

Synthesis Example 3: Synthesis of Compound 9

[0350]

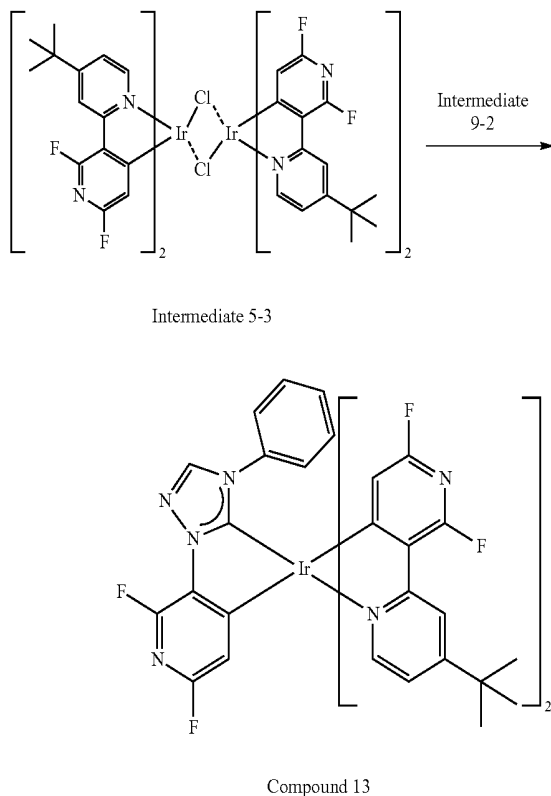


[0351] Intermediate 9-1 was obtained in the same manner as in the synthesis method of Intermediate 1-1 in Synthesis Example 1, except that iodobenzene was used instead of

iodomethane in synthesizing Intermediate 1-1, Intermediate 9-2 was obtained in the same manner as in the synthesis method of Intermediate 1-2 in Synthesis Example 1, except that Intermediate 9-1 was used instead of Intermediate 1-1 in synthesizing Intermediate 1-2, and 1.37 g (1.65 mmol, yield=15%) of Compound 9 was synthesized in the same manner as in Synthesis of Compound 1 in Synthesis Example 1, except that Intermediate 9-2 was used instead of Intermediate 1-2 in synthesizing Compound 1.

Synthesis Example 4: Synthesis of Compound 13

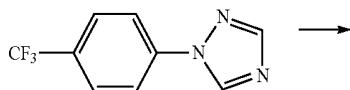
[0352]



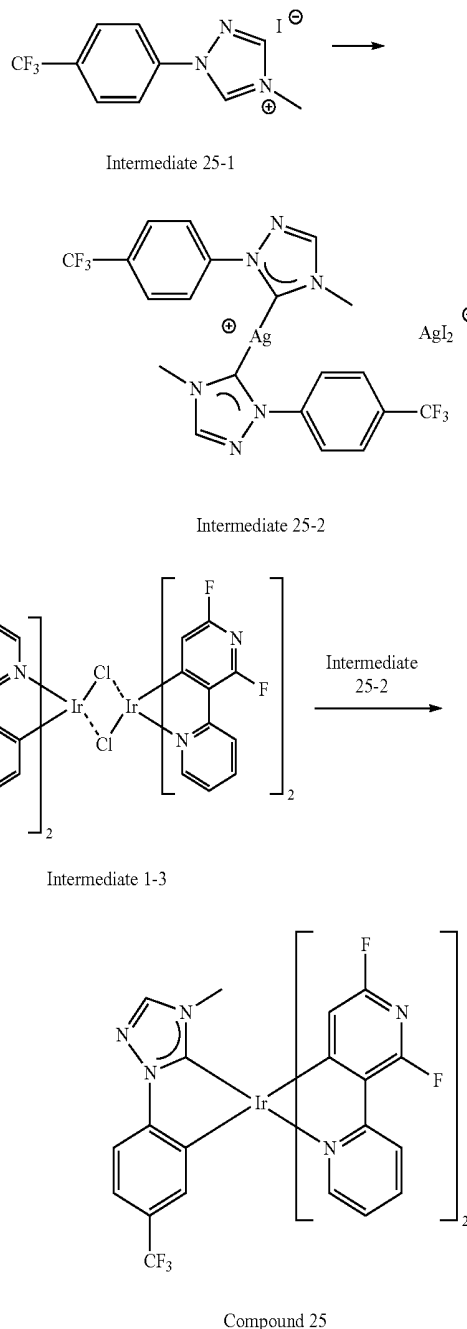
[0353] 1.56 g (1.65 mmol, yield=15%) of Compound 13 was synthesized in the same manner as in Synthesis of Compound 1 in Synthesis Example 1, except that Intermediates 9-2 and 5-3 are respectively used instead of Intermediates 1-2 and 1-3 in synthesizing Compound 1.

Synthesis Example 5: Synthesis of Compound 25

[0354]



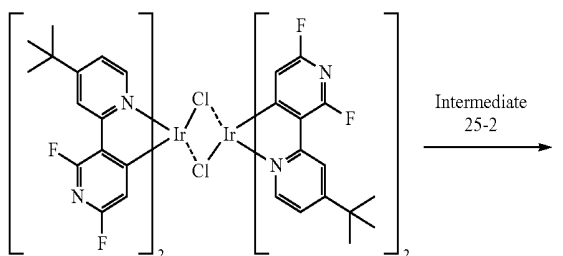
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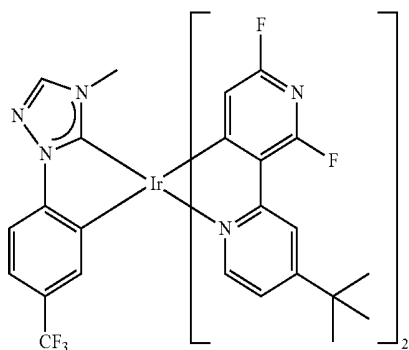
[0355] Intermediate 25-1 was obtained in the same manner as in the synthesis method of Intermediate 1-1 in Synthesis Example 1, except that 1-(4-(trifluoromethyl)phenyl)-1H-1,2,4-triazole was used instead of 2,6-difluoro-3-(1H-1,2,4-triazol-1-yl)pyridine in synthesizing Intermediate 1-1, Intermediate 25-2 was obtained in the same manner as in the synthesis method of Intermediate 1-2 in Synthesis Example 1, except that Intermediate 25-1 was used instead of Intermediate 1-1 in synthesizing Intermediate 1-2, and 1.58 g (1.98 mmol, yield=18%) of Compound 25 was synthesized in the same manner as in Synthesis of Compound 1 in Synthesis Example 1, except that Intermediate 25-2 was used instead of Intermediate 1-2 in synthesizing Compound 1.

Synthesis Example 6: Synthesis of Compound 29

[0356]



Intermediate 5-3

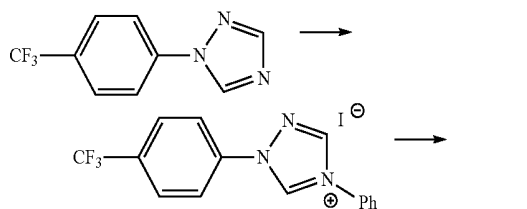


Compound 29

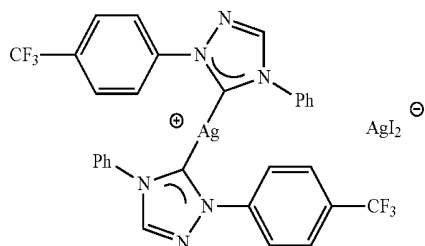
[0357] 1.61 g (1.76 mmol, yield=16%) of Compound 29 was synthesized in the same manner as in Synthesis of Compound 1 in Synthesis Example 1, except that Intermediates 25-2 and 5-3 were respectively used instead of Intermediates 1-2 and 1-3 in synthesizing Compound 1.

Synthesis Example 7: Synthesis of Compound 33

[0358]

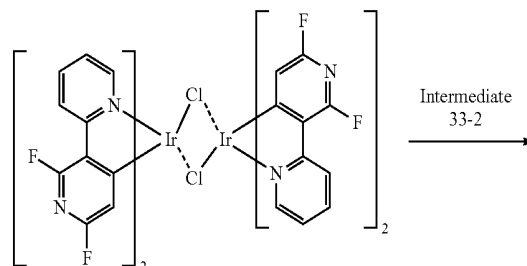


Intermediate 33-1

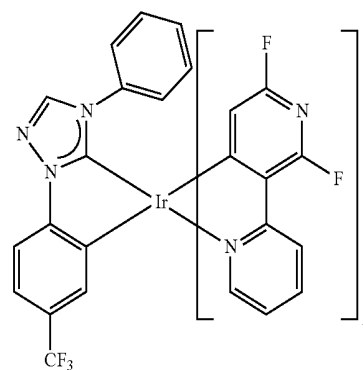


Intermediate 33-2

-continued



Intermediate 1-3

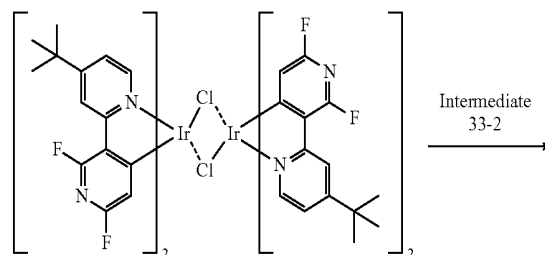


Compound 33

[0359] Intermediate 33-1 was obtained in the same manner as in the synthesis method of Intermediate 1-1 in Synthesis Example 1, except that 1-(4-(trifluoromethyl)phenyl)-1H-1,2,4-triazole and iodobenzene were respectively used instead of 2,6-difluoro-3-(1H-1,2,4-triazol-1-yl)pyridine and iodomethane in synthesizing Intermediate 1-1, Intermediate 33-2 was obtained in the same manner as in the synthesis method of Intermediate 1-2 in Synthesis Example 1, except that Intermediate 33-1 was used instead of Intermediate 1-1 in synthesizing Intermediate 1-2, and 1.56 g (1.65 mmol, yield=15%) of Compound 33 was synthesized in the same manner as in Synthesis of Compound 1 in Synthesis Example 1, except that Intermediate 33-2 was used instead of Intermediate 1-2 in synthesizing Compound 1.

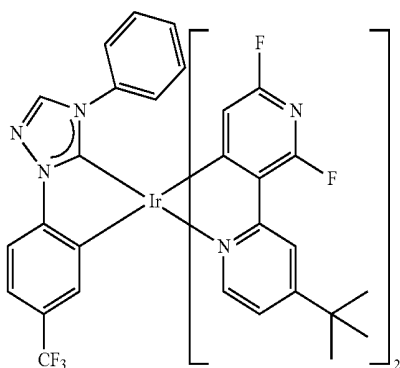
Synthesis Example 8: Synthesis of Compound 37

[0360]



Intermediate 5-3

-continued



Compound 37

[0361] 1.61 g (1.65 mmol, yield=15%) of Compound 37 was synthesized in Synthesis of Compound 1 in Synthesis Example 1, except that Intermediates 33-2 and 5-3 were respectively used instead of Intermediates 1-2 and 1-3 in synthesizing Compound 1.

Synthesis Example 9: Synthesis of Compound 49

[0362] 5.40 mmol of $[(CH_3)_2PtS(CH_3)]_2$, 2.70 mmol of Intermediate 1-2, and 2.70 mmol of 2',6'-difluoro-2,3'-bipyridine were mixed with 30 mL of tetrahydrofuran and then stirred at a temperature of 70° C. for 24 hours. The reaction mixture obtained therefrom was cooled to room temperature, and an extraction process was performed thereon by using dichloromethane and water. An organic layer obtained therefrom was dried by using magnesium sulfate and filtered by using celite, and column chromatography was performed thereon to obtain 1.92 g (3.29 mmol, yield=30%) of Compound 49.

Synthesis Example 10: Synthesis of Compound 2

[0363] 2.71 g (3.07 mmol, yield=28%) of Compound 53 was synthesized in the same manner as in Synthesis Example 9, except that 4-(tert-butyl)-2',6'-difluoro-2,3'-bipyridine was used instead of 2',6'-difluoro-2,3'-bipyridine.

[0364] ¹H NMR and FAB-MS of Compounds synthesized in Synthesis Examples 1 to 10 are shown in Table 1 below.

[0365] The synthesis methods of compounds other than Components shown in Table 1 may be easily recognized by those of ordinary skill in the art by referring to the synthesis mechanisms and source materials described above.

TABLE 1

Compound	¹ H NMR (CDCl ₃ , 500 MHz)	FAB-MS [M ⁺]	
		found	calc.
1	δ = 8.22 (dd, J = 13.5 Hz, J = 2.0 Hz, 2H), 7.98 (td, J = 8.4 Hz, J = 1.6 Hz, 2H), 7.89 (s, 1H), 7.86 (s, 1H), 7.60 (d, J = 7.5 Hz, 1H), 7.08 (t, J = 1.0 Hz, 1H), 6.93 (t, J = 1.0 Hz, 1H), 6.79 (d, J = 7.5 Hz, 1H), 6.10 (t, J = 2.0 Hz, 1H), 5.72 (t, J = 2.0 Hz, 1H), 3.29 (s, 3H).	770.16	770.10
5	δ = 8.23 (dd, J = 13.5 Hz, J = 2.0 Hz, 2H), 7.87 (s, 1H), 7.84 (s, 1H), 7.58 (d, J = 7.5 Hz, 1H), 7.10 (t, J = 1.0 Hz, 1H), 6.92 (t, J = 1.0 Hz, 1H), 6.77 (d, J = 7.5 Hz, 1H), 6.20 (t, J = 2.0 Hz, 1H), 5.52 (t, J = 2.0 Hz, 1H), 3.30 (s, 3H), 1.36 (s, 9H), 1.34 (s, 9H).	882.28	882.22
9	δ = 8.24 (dd, J = 13.5 Hz, J = 2.0 Hz, 2H), 7.93 (td, J = 8.4 Hz, J = 1.6 Hz, 2H), 7.91 (s, 1H), 7.88 (s, 1H), 7.62 (d, J = 7.5 Hz, 1H), 7.18 (t, J = 1.0 Hz, 1H), 6.92 (t, J = 1.0 Hz, 1H), 6.79 (d, J = 7.5 Hz, 1H), 6.72 (d, J = 6.5 Hz, 2H), 7.11 (td, J = 6.5 Hz, J = 2.1 Hz, 2H), 6.57-6.55 (m, 1H), 6.12 (t, J = 2.0 Hz, 1H), 5.70 (t, J = 2.0 Hz, 1H).	832.16	832.11
13	δ = 8.20 (dd, J = 13.5, J = 2.0 Hz, 2H), 7.90 (s, 1H), 7.87 (s, 1H), 7.65 (d, J = 7.5 Hz, 1H), 7.63 (td, J = 6.5 Hz, J = 2.1 Hz, 2H), 7.09 (t, J = 1.0 Hz, 1H), 6.56-6.54 (m, 1H), 6.92 (t, J = 1.0 Hz, 1H), 6.80 (d, J = 7.5 Hz, 1H), 6.77 (d, J = 6.5 Hz, 2H), 6.08 (t, J = 2.0 Hz, 1H), 5.70 (t, J = 2.0 Hz, 1H), 1.36 (s, 9H), 1.34 (s, 9H).	944.25	944.24
25	δ = 8.22 (dd, J = 13.5, J = 2.0 Hz, 2H), 7.96 (td, J = 8.4, J = 1.6 Hz, 2H), 7.93 (s, 1H), 7.85 (s, 1H), 7.72 (d, J = 8.0 Hz, 1H), 7.50 (d, J = 7.5 Hz, 1H), 7.41 (d, J = 8.0 Hz, 1H), 7.02 (t, J = 1.0 Hz, 1H), 6.93 (t, J = 1.0 Hz, 1H), 6.76 (d, J = 7.5 Hz, 1H), 6.09 (t, J = 2.0 Hz, 1H), 5.72 (t, J = 2.0 Hz, 1H), 3.28 (s, 3H).	801.15	801.11
29	δ = 8.20 (dd, J = 13.5, J = 2.0 Hz, 2H), 7.91 (s, 1H), 7.86 (s, 1H), 7.71 (d, J = 8.0 Hz, 1H), 7.61 (d, J = 7.5 Hz, 1H), 7.44 (d, J = 8.0 Hz, 1H), 7.08 (t, J = 1.0 Hz, 1H), 6.99 (t, J = 1.0 Hz, 1H), 6.81 (d, J = 7.5 Hz, 1H), 6.11 (t, J = 2.0 Hz, 1H), 5.73 (t, J = 2.0 Hz, 1H), 3.31 (s, 3H), 1.36 (s, 9H), 1.34 (s, 9H).	913.27	913.23

TABLE 1-continued

Compound	¹ H NMR (CDCl ₃ , 500 MHz)	FAB-MS [M ⁺]	
		found	calc.
33	δ = 8.21 (dd, J = 13.5, J = 2.0 Hz, 2H), 7.96 (td, J = 8.4, J = 1.6 Hz, 2H), 7.83 (s, 1H), 7.82 (s, 1H), 7.71 (d, J = 8.0 Hz, 1H), 7.62 (d, J = 7.5 Hz, 1H), 7.42 (d, J = 8.0 Hz, 1H), 7.39-7.22 (m, 5H), 7.10 (t, J = 1.0 Hz, 1H), 6.91 (t, J = 1.0 Hz, 1H), 6.77 (d, J = 7.5 Hz, 1H), 6.12 (t, J = 2.0 Hz, 1H), 5.71 (t, J = 2.0 Hz, 1H).	863.16	863.12
37	δ = 8.32 (dd, J = 13.5, J = 2.0 Hz, 2H), 7.91 (s, 1H), 7.86 (s, 1H), 7.74 (d, J = 8.0 Hz, 1H), 7.62 (d, J = 7.5 Hz, 1H), 7.44 (d, J = 8.0 Hz, 1H), 7.41-7.22 (m, 5H), 7.02 (t, J = 1.0 Hz, 1H), 6.94 (t, J = 1.0 Hz, 1H), 6.78 (d, J = 7.5 Hz, 1H), 6.11 (t, J = 2.0 Hz, 1H), 5.70 (t, J = 2.0 Hz, 1H), 1.35 (s, 9H), 1.34 (s, 9H).	975.28	975.25
49	δ = 8.52 (dd, pseudo-t, J = 25.2 Hz, 1H), 8.32 (d, J = 6.8 Hz, 1H), 8.04 (dd, pseudo-t, J = 24.9 Hz, 1H), 7.95 (s, 1H), 7.87 (s, 1H), 7.85-7.82 (m, 1H), 7.45-7.43 (m, 1H), 4.13 (s, 3H).	582.08	582.06
53	δ = 8.51 (dd, pseudo-t, J = 25.0 Hz, 1H), 8.30 (s, 1H), 8.02 (dd, pseudo-t, J = 24.9 Hz, 1H), 7.99 (s, 1H), 7.88 (s, 1H), 7.55 (d, J = 6.8 Hz, 1H), 4.12 (s, 3H), 1.32 (s, 9H).	638.55	638.51

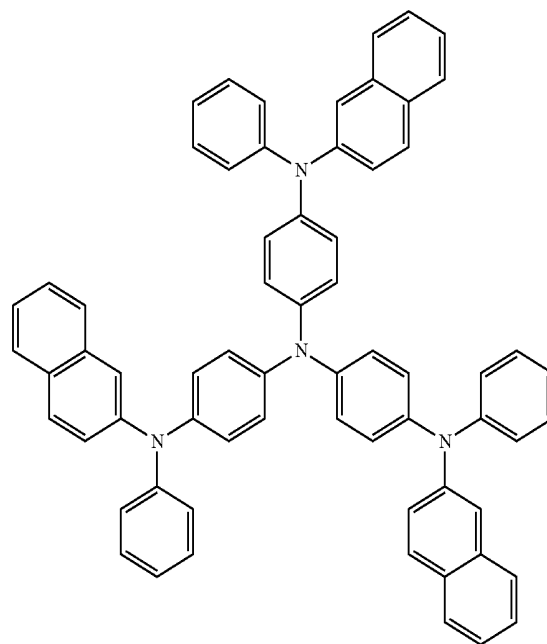
Example 1

[0366] An ITO anode was prepared by cutting a glass substrate (manufactured by Corning), on which an ITO layer was deposited to a thickness of 15 Ω/cm² (1,200 nm) and to a size of 50 mm×50 mm×0.7 mm, ultrasonically cleaning the glass substrate by using isopropyl alcohol, and pure water each for 5 minutes, and then, exposing the glass substrate to UV irradiation and ozone for 30 minutes to clean the ITO glass substrate. Then, the glass substrate was loaded into a vacuum deposition apparatus.

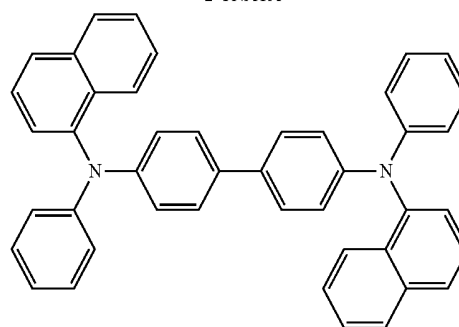
[0367] 2-TNATA was vacuum-deposited on the ITO anode formed on the glass substrate to form a hole injection layer having a thickness of 600 Å, and NPB was vacuum-deposited on the hole injection layer to form a hole transport layer having a thickness of 300 Å.

[0368] BCPDS and POPCPA (co-host) (a BCPDS to POPCPA weight ratio=1:1) and Compound 1 (dopant) were co-deposited on the hole transport layer at a co-host to dopant weight ratio of 90:10 to form an emission layer having a thickness of 300 Å.

[0369] TSPO1 was deposited on the emission layer to form a hole blocking layer having a thickness of 50 Å, Alq₃ was deposited on the hole blocking layer to form an electron transport layer having a thickness of 300 Å, LiF was deposited on the electron transport layer to form an electron injection layer having a thickness of 10 Å, and Al was vacuum-deposited on the electron injection layer to form a cathode having a thickness of 3,000 Å, thereby completing the manufacture of an organic light-emitting device.

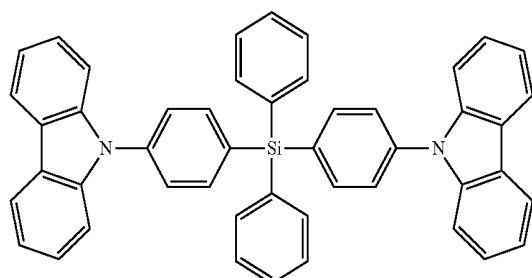


2-TNATA



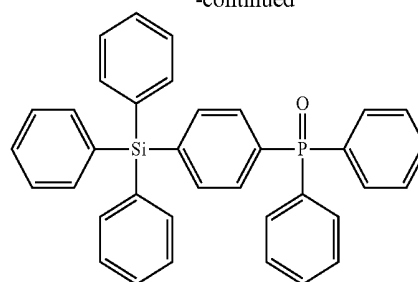
NPB

-continued



BCPDS

-continued



TSP01

Examples 2 to 10 and Comparative Examples A to C

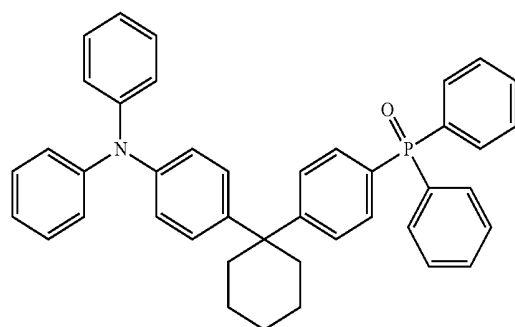
[0370] Organic light-emitting devices of Examples 2 to 10 and Comparative Examples A to C were manufactured in the same manner as in Example 1, except that Compounds shown in Table 2 were each used instead of Compound 1 as a dopant in forming an emission layer.

Evaluation Example 1

[0371] The driving voltage, current density, luminance, luminescent efficiency, and maximum emission wavelength of the organic light-emitting devices manufactured according to Examples 1 to 10 and Comparative Examples A to C were measured by using a Keithley SMU 236 and a luminance meter PR650. Results thereof are shown in Table 2 below.

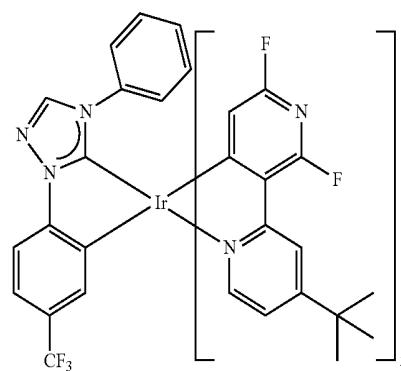
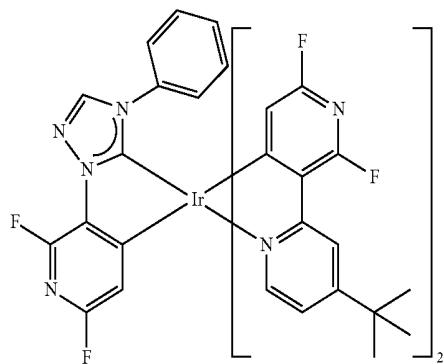
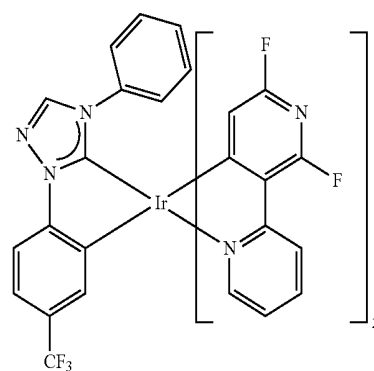
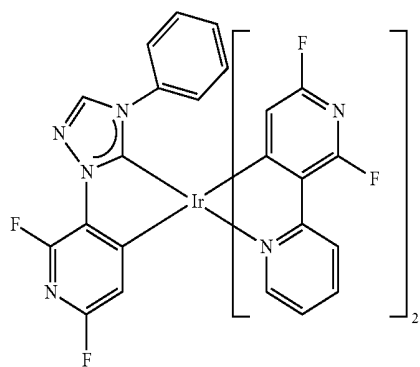
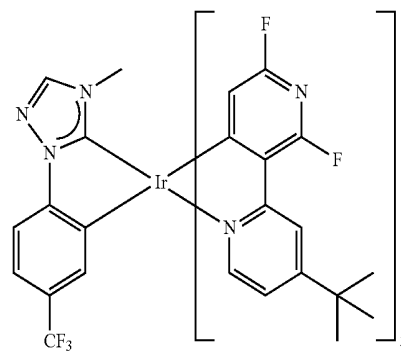
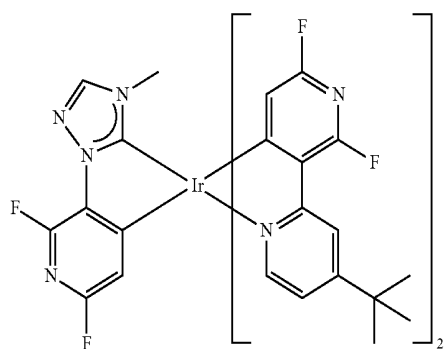
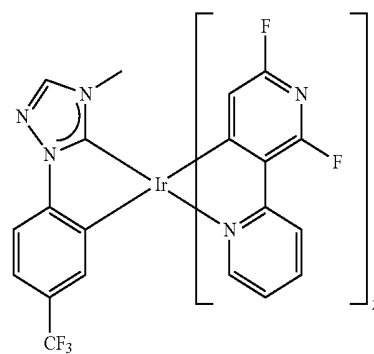
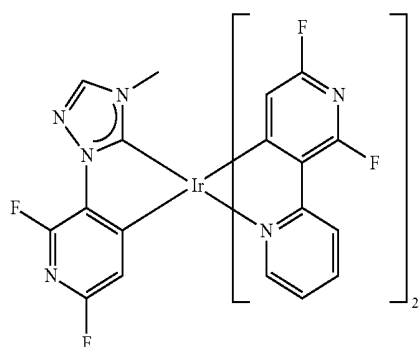
TABLE 2

Dopant compound No.	Driving voltage (V)	Current density (mA/cm ²)	Luminance (cd/m ²)	Luminescent Efficiency (cd/A)	Emission color	Maximum emission wavelength (nm)
Example 1	1	3.4	0.09	15	Blue	449
Example 2	5	3.4	0.08	15	Blue	449
Example 3	9	3.3	0.08	15	Blue	452
Example 4	13	3.3	0.08	15	Blue	452
Example 5	25	3.3	0.09	15	Blue	450
Example 6	29	3.3	0.08	15	Blue	451
Example 7	33	3.4	0.08	15	Blue	455
Example 8	37	3.4	0.06	15	Blue	456
Example 9	49	3.3	0.10	15	Blue	448
Example 10	53	3.4	0.10	15	Blue	448
Comparative Example A	A	4.3	0.25	15	Sky blue	483
Comparative Example B	B	4.2	0.18	15	Sky blue	475
Comparative Example C	C	3.9	0.17	15	Blue	468

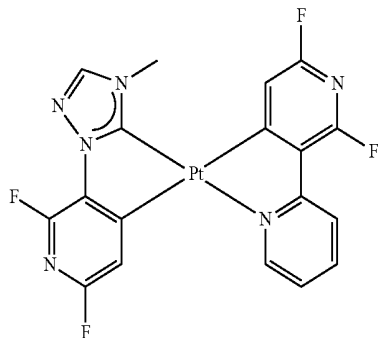


POPCPA

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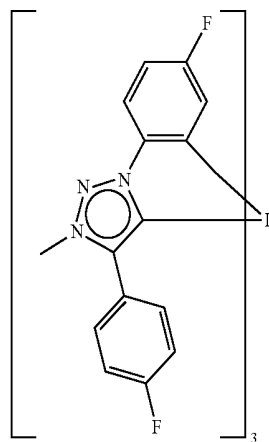


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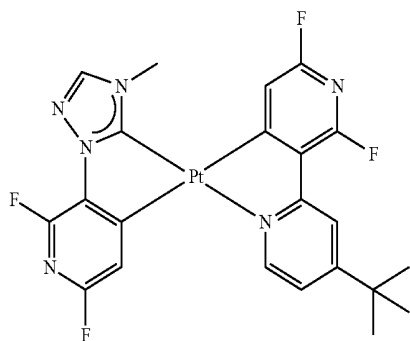
-continued

49



<Compound C>

53



[0372] Referring to Table 2, it can be confirmed that the organic light-emitting devices of Examples 1 to 10 have a lower driving voltage, and a higher luminescent efficiency and are suitable for emission of deep blue light, as compared with those of the organic light-emitting devices of Comparative Examples A to C.

[0373] According to one or more embodiments, an organic light-emitting device including the organometallic compound may have a low driving voltage, and high luminescent efficiency.

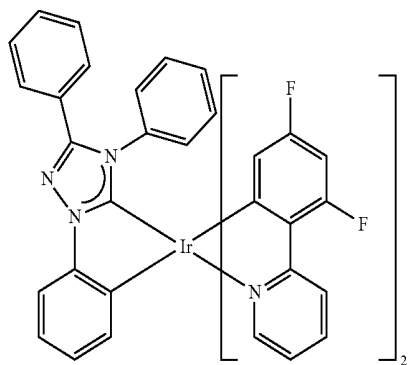
[0374] It should be understood that embodiments described herein should be considered in a descriptive sense only and not for purposes of limitation. Descriptions of features or aspects within each embodiment should typically be considered as available for other similar features or aspects in other embodiments.

[0375] While one or more embodiments have been described with reference to the figures, it will be understood by those of ordinary skill in the art that various changes in form and details may be made therein without departing from the spirit and scope as defined by the following claims.

What is claimed is:

1. An organometallic compound represented by Formula 1:

<Compound A>

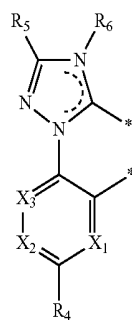
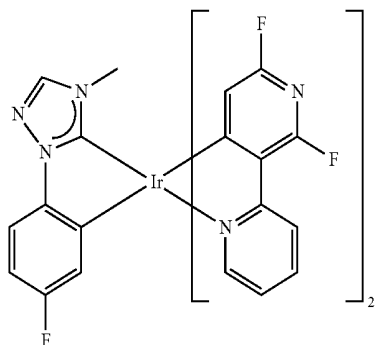


<Formula 1>

$M(L_1)_{n1}(L_2)_{n2}$

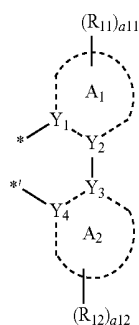
<Formula 2A>

<Compound B>



-continued

<Formula 2B>



wherein M in Formula 1 is a third row transition metal, in Formula 1, L_1 is a ligand represented by Formula 2A, and n_2 is 1, 2, or 3, wherein, when n_2 is two or more, two or more $L_1(s)$ are identical to or different from each other,

in Formula 1, L_2 is a ligand represented by Formula 2B, and n_2 is 0, 1, or 2, wherein, when n_2 is two or more, two or more $L_2(s)$ are identical to or different from each other,

the sum of n_1 and n_2 in Formula 1 is two or three,

* and * in Formulae 2A and 2B each indicate a binding site to M in Formula 1,

in Formula 2A, X_1 is N or C(R_1), X_2 is N or C(R_2), and X_3 is N or C(R_3),

in Formula 2B, Y_1 is carbon, Y_2 and Y_3 are each independently carbon or nitrogen, Y_4 is nitrogen, a bond between Y_1 and Y_2 is a single bond or a double bond, a bond between Y_2 and Y_3 is a single bond, and a bond between Y_3 and Y_4 is a single bond or a double bond,

ring A_1 and ring A_2 in Formula 2B are each a C_2 - C_{60} heterocyclic group comprising at least one N as a ring-forming atom,

in Formulae 2A and 2B, R_1 to R_3 , R_5 , R_6 , R_{11} , and R_{12} are each independently selected from hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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_3 - C_{10} cycloalkenyl 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), —N(Q_1)(Q_2), —B(Q_1)(Q_2), —C(=O)(Q_1), —S(=O)₂(Q_1), and —P(=O)(Q_1)(Q_2),

R_4 is an electron withdrawing group,

R_2 and R_3 in Formula 2A are optionally linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group,

R_5 and R_6 in Formula 2A are optionally linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group,

two neighboring groups of a plurality of $R_{11}(s)$ in Formula 2B are optionally linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group,

two neighboring groups of a plurality of $R_{12}(s)$ in Formula 2B are optionally linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group,

neighboring R_{11} and R_{12} in Formula 2B are optionally linked to form a substituted or unsubstituted C_4 - C_{60} carbocyclic group or a substituted or unsubstituted C_2 - C_{60} heterocyclic group,

a_{11} and a_{12} in Formula 2B are each independently an integer from 0 to 10, and

at least one substituent of the substituted C_1 - C_{60} alkyl group, the substituted C_2 - C_{60} alkenyl group, the substituted C_2 - C_{60} alkynyl group, the substituted C_2 - C_{60} alkoxy group, the substituted C_3 - C_{10} cycloalkyl group, the substituted C_1 - C_{10} heterocycloalkyl group, the substituted C_3 - C_{10} cycloalkenyl group, the substituted C_1 - C_{10} heterocycloalkenyl group, the substituted C_6 - C_{60} aryl group, the substituted C_6 - C_{60} aryloxy group, the substituted C_6 - C_{60} arylthio group, the substituted C_1 - C_{60} heteroaryl group, the substituted monovalent non-aromatic condensed polycyclic group, the substituted monovalent non-aromatic condensed heteropolycyclic group, the substituted C_4 - C_{60} carbocyclic group, and the substituted C_2 - C_{60} heterocyclic group is selected from:

deuterium (-D), —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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;

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 deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C_3 - C_{10} cycloalkyl group, a C_1 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_1 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, a C_1 - C_{60} heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, —Si(Q_{11})(Q_{12})(Q_{13}), —N(Q_{11})(Q_{12}), —B(Q_{11})(Q_{12}), —C(=O)(Q_{11}), —S(=O)₂(Q_{11}), and —P(=O)(Q_{11})(Q_{12});

a C_3 - C_{10} cycloalkyl group, a C_1 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_1 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, a C_1 - C_{60} heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic 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 deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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), —N(Q₂₁)(Q₂₂), —B(Q₂₁)(Q₂₂), —C(=O)(Q₂₁), —S(=O)₂(Q₂₁), and —P(=O)(Q₂₁)(Q₂₂); and —Si(Q₃₁)(Q₃₂)(Q₃₃), —N(Q₃₁)(Q₃₂), —B(Q₃₁)(Q₃₂), —C(=O)(Q₃₁), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂), and Q₁ to Q₃, Q₁₁ to Q₁₃, Q₂₁ to Q₂₃, and Q₃₁ to Q₃₃ are each independently selected from hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, 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, a monovalent non-aromatic condensed heteropolycyclic group, a C₁-C₆₀ alkyl group substituted with at least one selected from deuterium, —F, and a cyano group, a C₆-C₆₀ aryl group substituted with at least one selected from deuterium, —F, and a cyano group, a biphenyl group, and a terphenyl group.
2. The organometallic compound of claim 1, wherein M in Formula 1 is iridium or osmium and the sum of n1 and n2 is three; or M in Formula 1 is platinum and the sum of n1 and n2 is two.
3. The organometallic compound of claim 1, wherein in Formula 2A, X₁ is C(R₁), X₂ is C(R₂), and X₃ is C(R₃); or in Formula 2A, X₁ is C(R₁), X₂ is N, and X₃ is C(R₃).
4. The organometallic compound of claim 1, wherein, in Formula 2A, X₃ is C(R₃) and R₃ is an electron withdrawing group.
5. The organometallic compound of claim 1, wherein ring A₁ and ring A₂ in Formula 2B are each independently a pyridine group, a pyrimidine group, a pyrazine group, a pyridazine group, a triazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, a phenanthroline group, a pyrazole group, an imidazole group, a triazole group, an oxazole group, an iso-oxazole group, a thiazole group, an isothiazole group, an oxadiazole group, a thiadiazol group, a benzopyrazole group, a benzimidazole group, a benzoxazole group, a benzothiazole group, a benzoxadiazole group, a benzothiadiazol group, a 5,6,7,8-tetrahydroisoquinoline group, or a 5,6,7,8-tetrahydroquinoline group.
6. The organometallic compound of claim 1, wherein R₁ to R₃, R₅, R₆, R₁₁, and R₁₂ in Formulae 2A and 2B are each independently selected from: hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, and a C₁-C₂₀ alkoxy group; a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, and a C₁-C₂₀ alkoxy group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, —CD₃, —CD₂H, —CDH₂, —CF₃, —CF₂H, —CFH₂, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a pyridinyl group, a pyrimidinyl group, —Si(Q₃₁)(Q₃₂)(Q₃₃), —N(Q₃₁)(Q₃₂), —B(Q₃₁)(Q₃₂), —C(=O)(Q₃₁), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂); a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl group, a benzofuranlyl 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 dibenzofuranlyl group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, an imidazopyridinyl group, and an imidazopyrimidinyl group; a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl

group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazoliny group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl 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, an imidazopyridinyl group, and an imidazopyrimidinyl group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, —CD₃, —CD₂H, —CDH₂, —CF₃, —CF₂H, —CFH₂, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, a C₁-C₂₀ alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazoliny group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl 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, an imidazopyridinyl group, an imidazopyrimidinyl group, —Si(Q₃₁)(Q₃₂)(Q₃₃), —N(Q₃₁)(Q₃₂), —B(Q₃₁)(Q₃₂), —C(=O)(Q₃₁), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂); and —Si(Q₁)(Q₂)(Q₃), —N(Q₁)(Q₂), —B(Q₁)(Q₂), —C(=O)(Q₁), —S(=O)₂(Q₁), and —P(=O)(Q₁)(Q₂), and

Q₁ to Q₃ and Q₃₁ to Q₃₃ are each independently selected from:

hydrogen, deuterium, —F, —Cl, —Br, —I, a cyano group, a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, a C₁-C₂₀ alkoxy group, a C₃-C₁₀ cycloalkyl group, a C₁-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₁-C₁₀ heterocycloalkenyl group, a C₆-C₂₀ aryl group, a C₁-C₂₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group;

a C₁-C₂₀ alkyl group substituted with at least one selected from deuterium, —F, and a cyano group;

a C₁-C₂₀ aryl group substituted with at least one selected from deuterium, —F, and a cyano group; and

a biphenyl group and a terphenyl group.

7. The organometallic compound of claim 1, wherein, in Formula 2B, the sum of a11 and a12 is equal to or greater than one, and at least one selected from R₁₁ and R₁₂ is an electron withdrawing group.

8. The organometallic compound of claim 1, wherein the electron withdrawing group is —F, —Cl, —Br, —I, a cyano group, a nitro group, or —B(Q₁)(Q₂); or a first group substituted with at least one selected from —F, —Cl, —Br, —I, —CF₃, —CF₂H, —CFH₂, a cyano group, a nitro group, and —B(Q₃₁)(Q₃₂),

provided that the first group is selected from:

a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, and a C₁-C₂₀ alkoxy group;

a C₁-C₂₀ alkyl group, a C₂-C₂₀ alkenyl group, a C₂-C₂₀ alkynyl group, and a C₁-C₂₀ alkoxy group, each substituted with at least one selected from deuterium, —CD₃, —CD₂H, —CDH₂, a hydroxyl group, an amidino group, a hydrazino group, a hydrazono group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a pyridinyl group, a pyrimidinyl group, —Si(Q₃₁)(Q₃₂)(Q₃₃), —N(Q₃₁)(Q₃₂), —C(=O)(Q₃₁), —S(=O)₂(Q₃₁), and —P(=O)(Q₃₁)(Q₃₂);

a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazoliny group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl 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, an imidazopyridinyl group, and an imidazopyrimidinyl group; and

a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl

group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazoliny group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl 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, an imidazopyridinyl group, and an imidazopyrimidinyl group, each substituted with at least one selected from deuterium, $-\text{CD}_3$, $-\text{CD}_2\text{H}$, $-\text{CDH}_2$, a hydroxyl group, an amidino group, a hydrazino group, a hydrazono group, a $\text{C}_1\text{-C}_{20}$ alkyl group, a $\text{C}_2\text{-C}_{20}$ alkenyl group, a $\text{C}_2\text{-C}_{20}$ alkynyl group, a $\text{C}_1\text{-C}_{20}$ alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a fluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl 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, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a quinoxalinyl group, a quinazoliny group, a cinnolinyl group, a carbazolyl group, a phenanthrolinyl group, a benzimidazolyl 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, an imidazopyridinyl group, an imidazopyrimidinyl group, $-\text{Si}(\text{Q}_{31})(\text{Q}_{32})(\text{Q}_{33})$, $-\text{N}(\text{Q}_{31})(\text{Q}_{32})$, $-\text{C}(=\text{O})(\text{Q}_{31})$, $-\text{S}(=\text{O})_2(\text{Q}_{31})$, and $-\text{P}(=\text{O})(\text{Q}_{31})(\text{Q}_{32})$, and

Q_1 to Q_3 and Q_{31} to Q_{33} are each independently selected from:

hydrogen, deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a cyano group, a $\text{C}_1\text{-C}_{20}$ alkyl group, a $\text{C}_2\text{-C}_{20}$ alkenyl group, a $\text{C}_2\text{-C}_{20}$ alkynyl group, a $\text{C}_1\text{-C}_{20}$ alkoxy group, 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 $\text{C}_6\text{-C}_{20}$ aryl group, a $\text{C}_1\text{-C}_{20}$ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group;

a $\text{C}_1\text{-C}_{20}$ alkyl group substituted with at least one selected from deuterium, $-\text{F}$, and a cyano group;

a $\text{C}_6\text{-C}_{20}$ aryl group substituted with at least one selected from deuterium, $-\text{F}$, and a cyano group; and

a biphenyl group and a terphenyl group.

9. The organometallic compound of claim 1, wherein

the electron withdrawing group is $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a cyano group, a nitro group, or $-\text{B}(\text{Q}_1)(\text{Q}_2)$; or a first group substituted with at least one selected from $-\text{F}$,

$-\text{Cl}$, $-\text{Br}$, $-\text{I}$, $-\text{CF}_3$, $-\text{CF}_2\text{H}$, $-\text{CFH}_2$, a cyano group, a nitro group, and $-\text{B}(\text{Q}_{31})(\text{Q}_2)$,

provided that the first group is selected from:

a $\text{C}_1\text{-C}_{20}$ alkyl group, a $\text{C}_2\text{-C}_{20}$ alkenyl group, a $\text{C}_2\text{-C}_{20}$ alkynyl group, and a $\text{C}_1\text{-C}_{20}$ alkoxy group;

a $\text{C}_1\text{-C}_{20}$ alkyl group, a $\text{C}_2\text{-C}_{20}$ alkenyl group, a $\text{C}_2\text{-C}_{20}$ alkynyl group, and a $\text{C}_1\text{-C}_{20}$ alkoxy group, each substituted with at least one selected from deuterium, $-\text{CD}_3$, $-\text{CD}_2\text{H}$, $-\text{CDH}_2$, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, a pyridinyl group, and a pyrimidinyl group;

a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, and a fluorenyl group; and

a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, and a fluorenyl group, each substituted with at least one selected from deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, $-\text{CD}_3$, $-\text{CD}_2\text{H}$, $-\text{CDH}_2$, a $\text{C}_1\text{-C}_{20}$ alkyl group, a $\text{C}_2\text{-C}_{20}$ alkenyl group, a $\text{C}_2\text{-C}_{20}$ alkynyl group, a $\text{C}_1\text{-C}_{20}$ alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclooctyl group, an adamantanyl group, a norbornanyl group, a norbornenyl group, a cyclopentenyl group, a cyclohexenyl group, a cycloheptenyl group, a phenyl group, a naphthyl group, and a fluorenyl group, and

Q_1 , Q_2 , Q_{31} , and Q_{32} are each independently selected from:

hydrogen, deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a cyano group, a $\text{C}_1\text{-C}_{20}$ alkyl group, a $\text{C}_1\text{-C}_{20}$ alkoxy group, a phenyl group, and a naphthyl group;

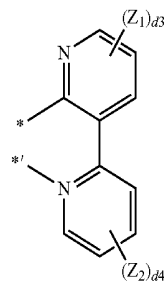
a $\text{C}_1\text{-C}_{20}$ alkyl group substituted with at least one selected from deuterium, $-\text{F}$, and a cyano group;

a phenyl group and a naphthyl group, each substituted with at least one selected from deuterium, $-\text{F}$, and a cyano group; and

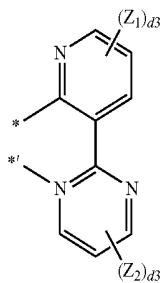
a biphenyl group and a terphenyl group.

10. The organometallic compound of claim 1, wherein L_2 is a ligand selected from groups represented by Formulae 2B-1 to 2B-60:

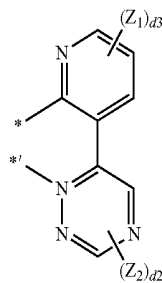
Formula 2B-1



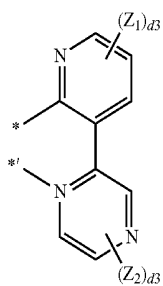
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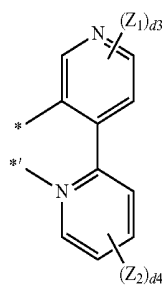
Formula 2B-2



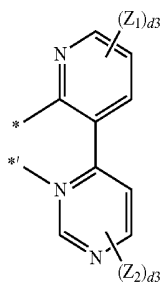
Formula 2B-7



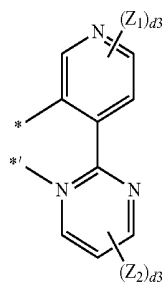
Formula 2B-3



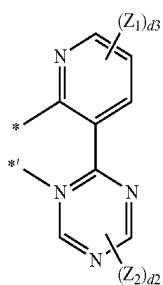
Formula 2B-8



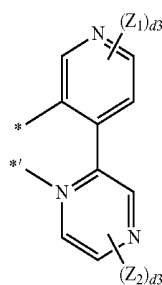
Formula 2B-4



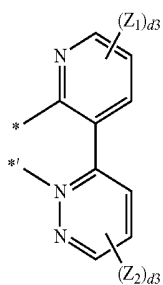
Formula 2B-9



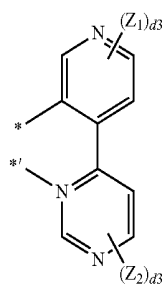
Formula 2B-5



Formula 2B-10

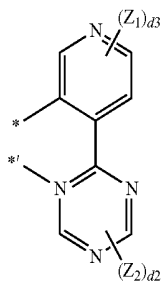


Formula 2B-6

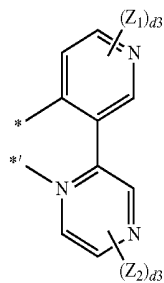


Formula 2B-11

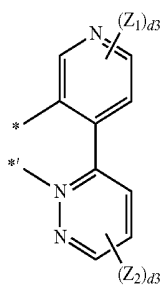
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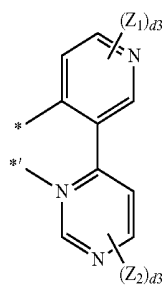
Formula 2B-12



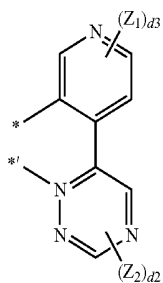
Formula 2B-17



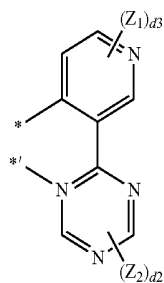
Formula 2B-13



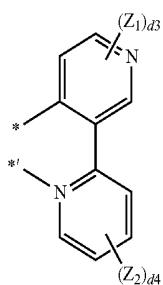
Formula 2B-18



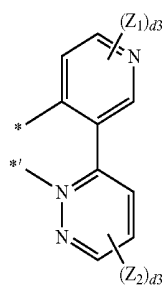
Formula 2B-14



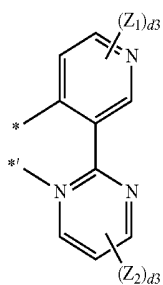
Formula 2B-19



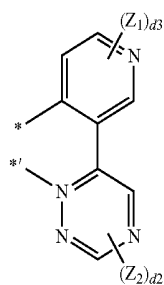
Formula 2B-15



Formula 2B-20

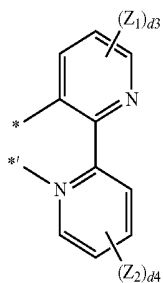


Formula 2B-16



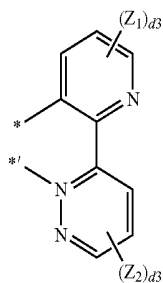
Formula 2B-21

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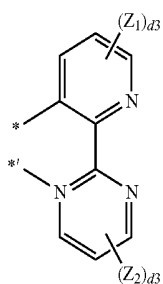


Formula 2B-22

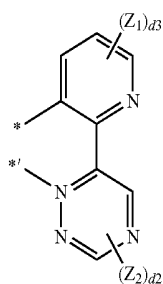
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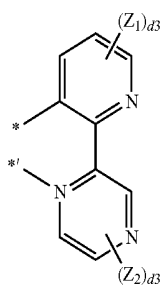
Formula 2B-27



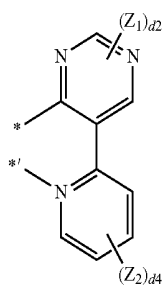
Formula 2B-23



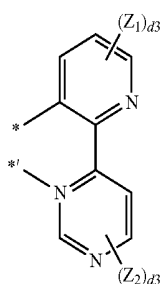
Formula 2B-28



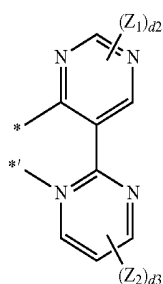
Formula 2B-24



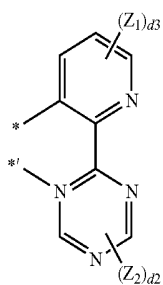
Formula 2B-29



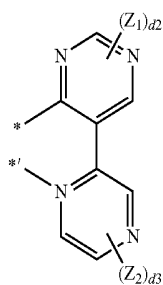
Formula 2B-25



Formula 2B-30

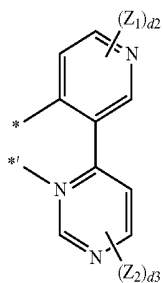


Formula 2B-26

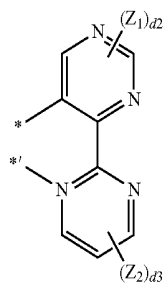


Formula 2B-31

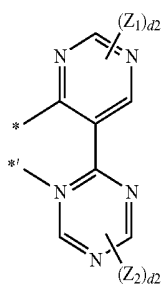
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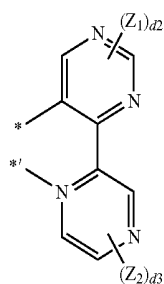
Formula 2B-32



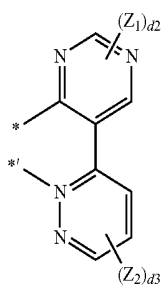
Formula 2B-37



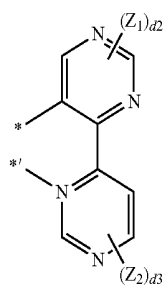
Formula 2B-33



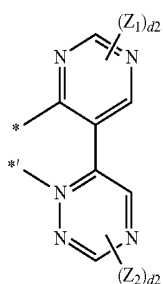
Formula 2B-38



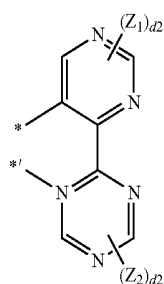
Formula 2B-34



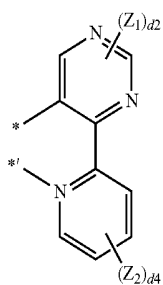
Formula 2B-39



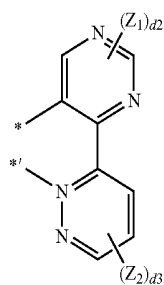
Formula 2B-35



Formula 2B-40

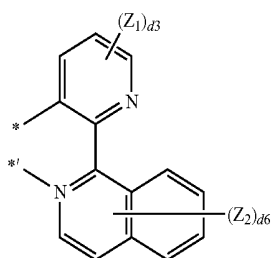


Formula 2B-36

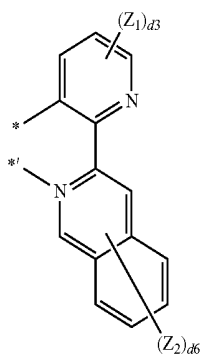


Formula 2B-41

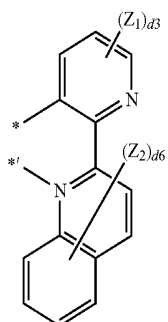
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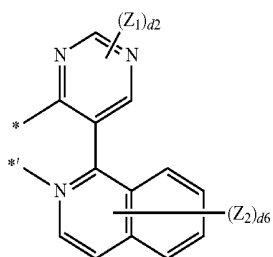
Formula 2B-52



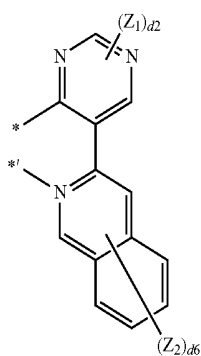
Formula 2B-53



Formula 2B-54

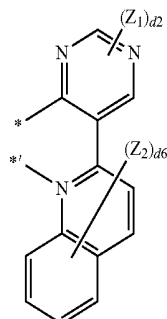


Formula 2B-55

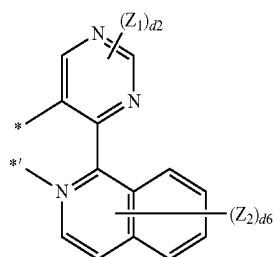


Formula 2B-56

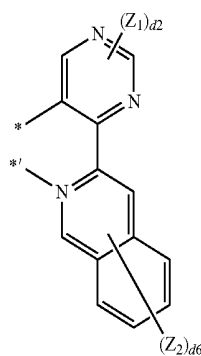
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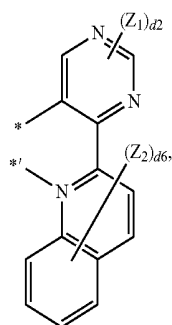
Formula 2B-57



Formula 2B-58



Formula 2B-59



Formula 2B-60

wherein, in Formulae 2B-1 to 2B-60,

Z_1 and Z_2 are the same as described in connection with R_{11} and R_{12} in claim 1, respectively,

d_2 is an integer from 0 to 2,

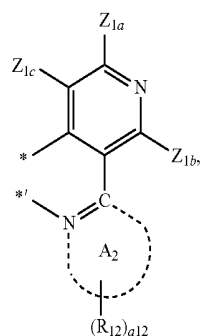
d_3 is an integer from 0 to 3,

d_4 is an integer from 0 to 4,

d_6 is an integer from 0 to 6, and

* and *' each indicate a binding site to M in Formula 1.

11. The organometallic compound of claim 1, wherein L_2 is a ligand selected from groups represented by Formula 2B-A:



<Formula 2B-A>

wherein, in Formula 2B-A,

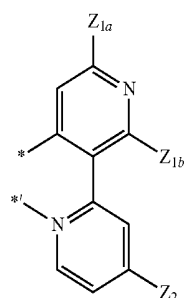
Z_{1a} , Z_{1b} , and Z_{1c} are the same as described in connection with R_{11} in claim 1,

ring A_2 is a pyridine group, a pyrimidine group, a pyrazine group, a pyridazine group, a triazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, a phenanthroline group, a 5,6,7,8-tetrahydroisoquinoline group, or a 5,6,7,8-tetrahydroquinoline group,

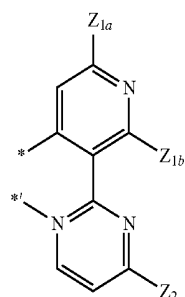
R_{12} and a_{12} are the same as described in claim 1, and * and *' each indicate a binding site to M in Formula 1.

12. The organometallic compound of claim 11, wherein at least one selected from Z_{1a} and Z_{1b} is an electron withdrawing group.

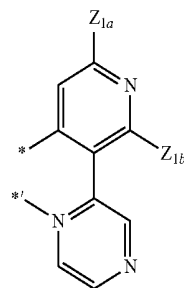
13. The organometallic compound of claim 1, wherein L_2 is a ligand selected from groups represented by Formulae 2B(1) to 2B(8):



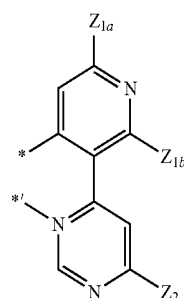
Formula 2B(1)



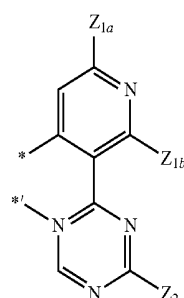
Formula 2B(2)



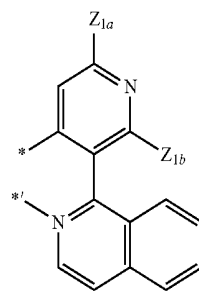
Formula 2B(3)



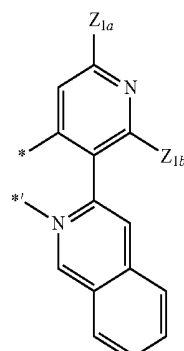
Formula 2B(4)



Formula 2B(5)

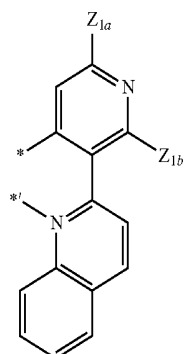


Formula 2B(6)



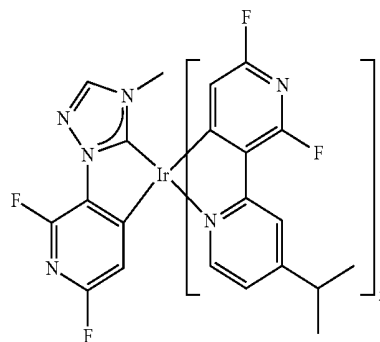
Formula 2B(7)

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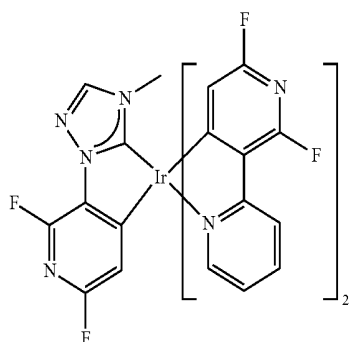
Formula 2B(8)

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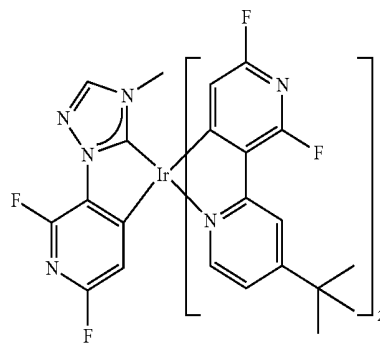


4

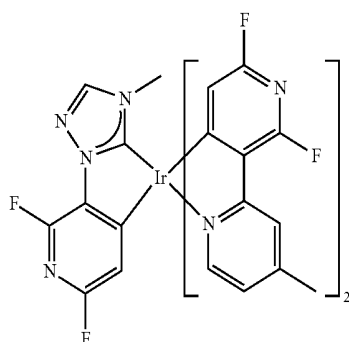
wherein, in Formulae 2B(1) and 2B(8),
 Z_{1a} and Z_{1b} are the same as described in connection with R_{11} in claim 1,
 Z_2 is the same as described with R_{12} in claim 1, and
 * and *' each indicate a binding site to M in Formula 1.
14. The organometallic compound of claim 1, wherein the organometallic compound is selected from Compounds 1 to 96:



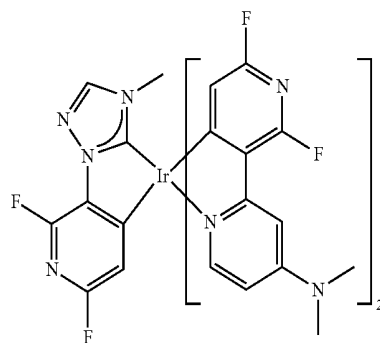
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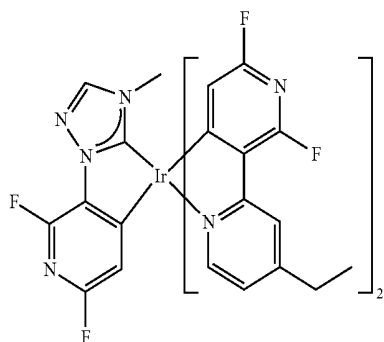
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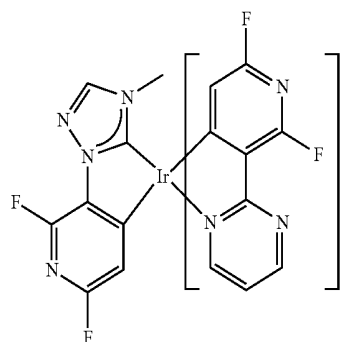
2



6

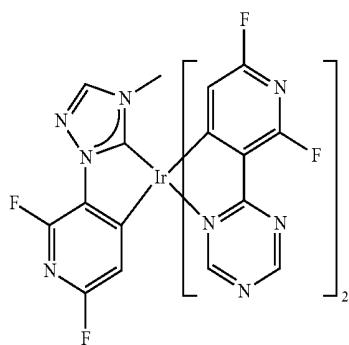


3



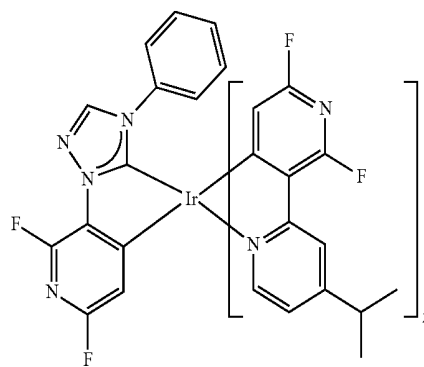
7

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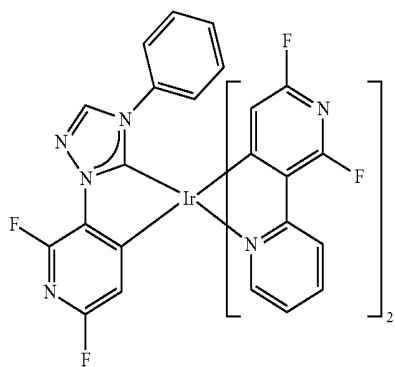
8

12



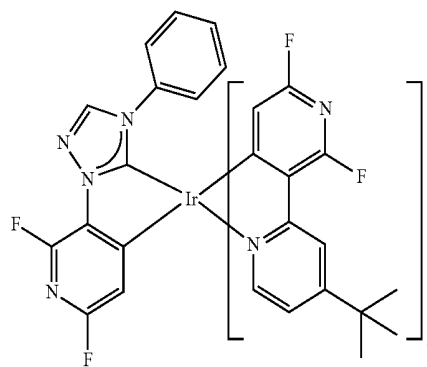
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13



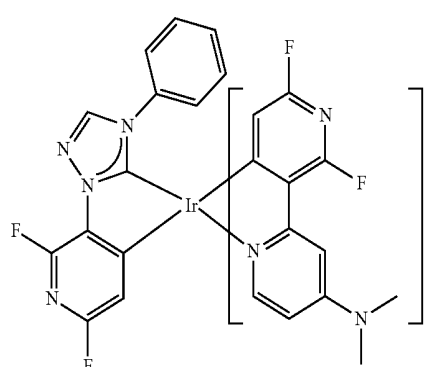
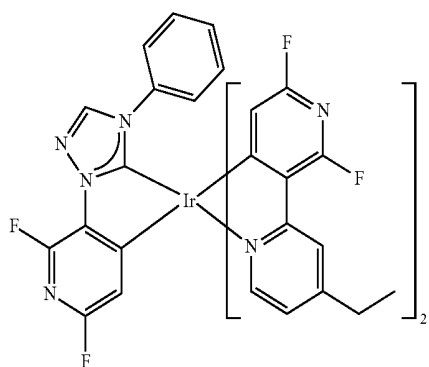
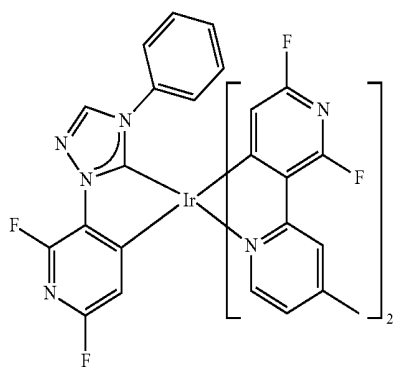
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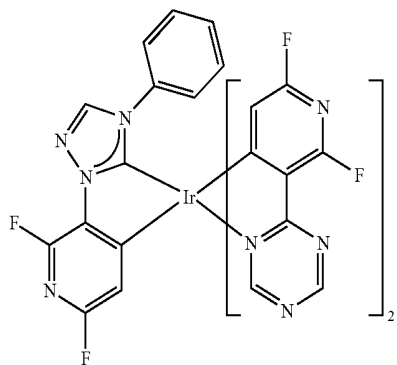


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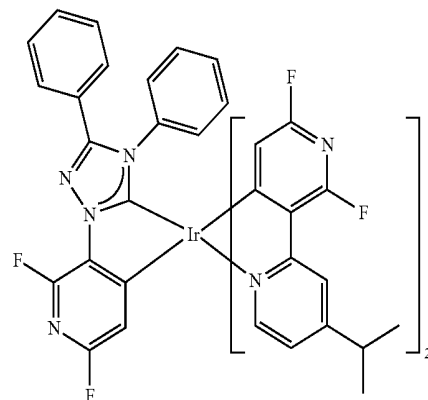


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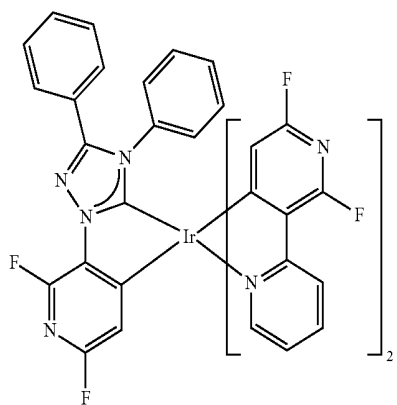
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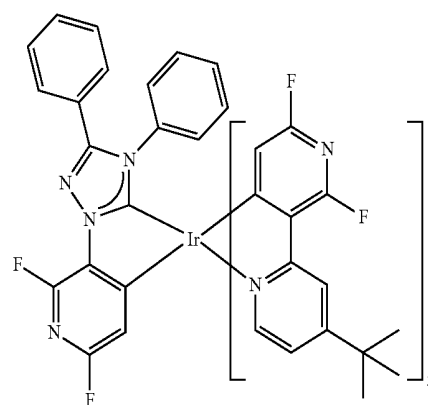
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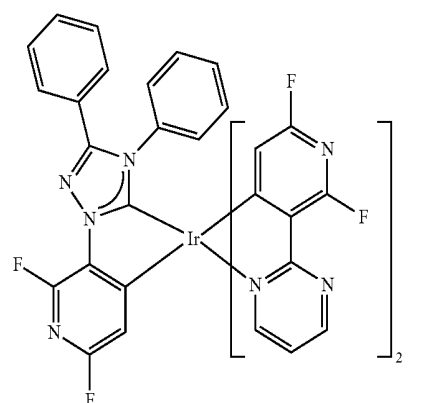
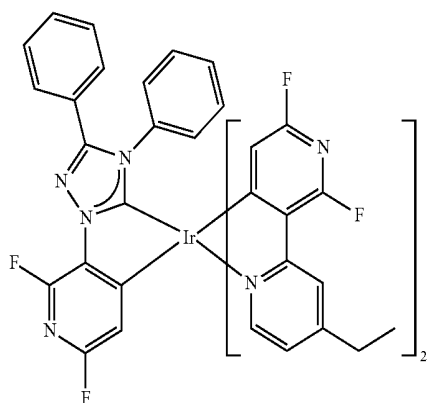
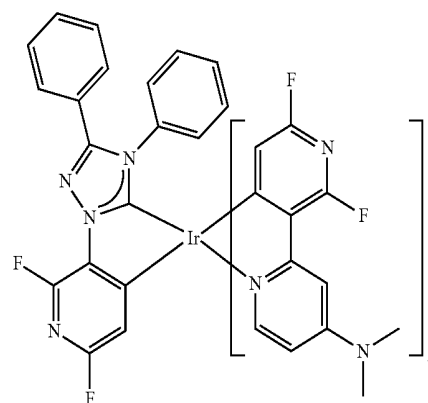
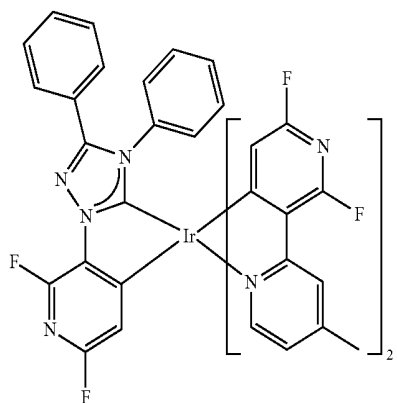
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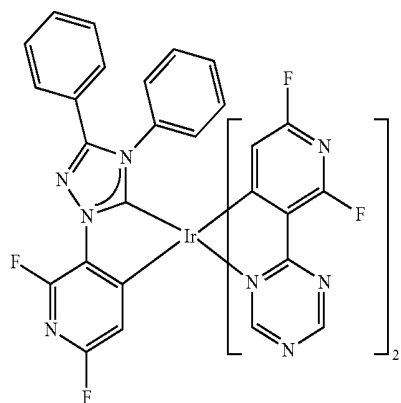


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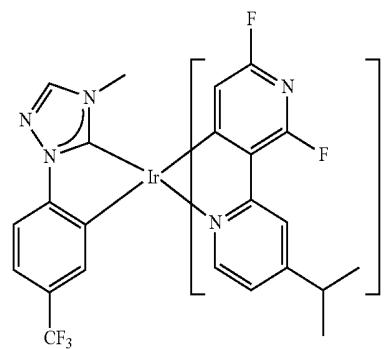


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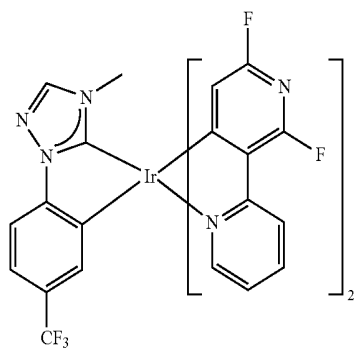


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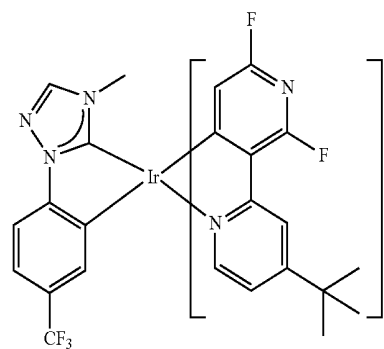
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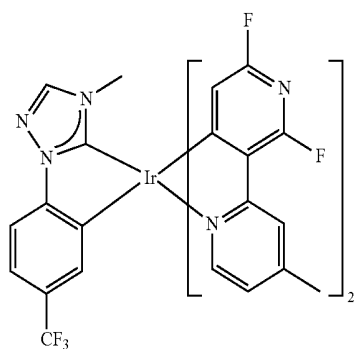
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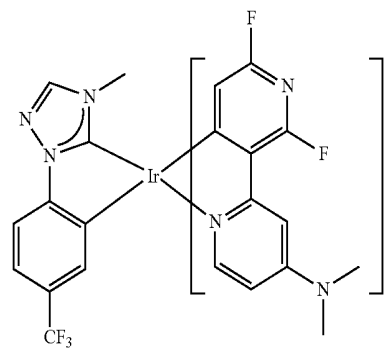
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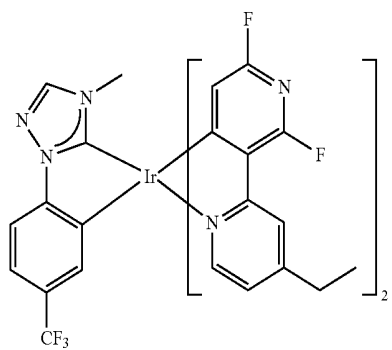
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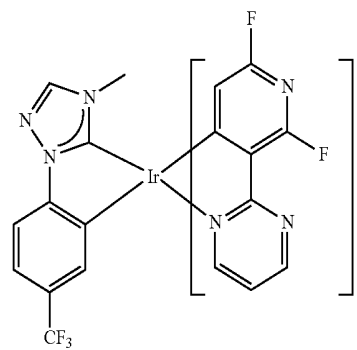
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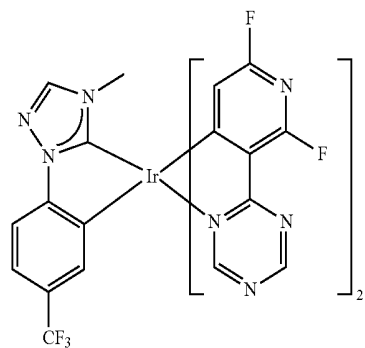


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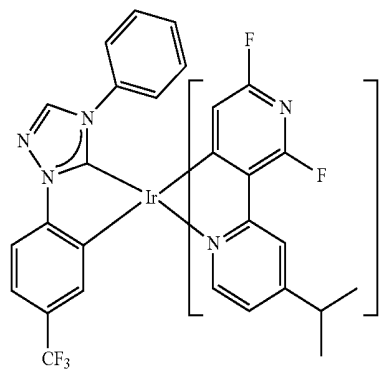
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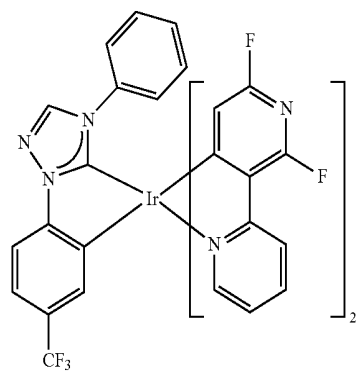


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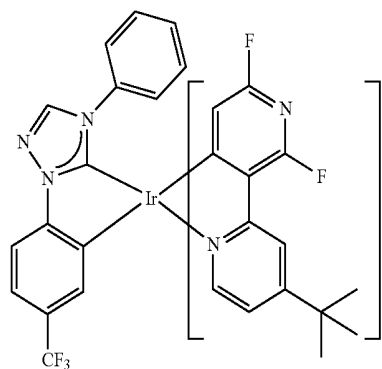
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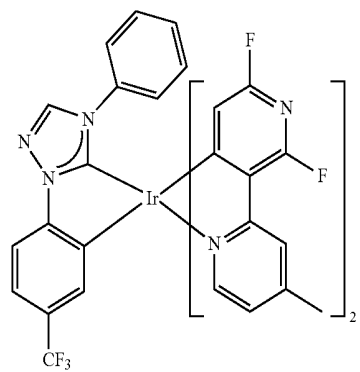
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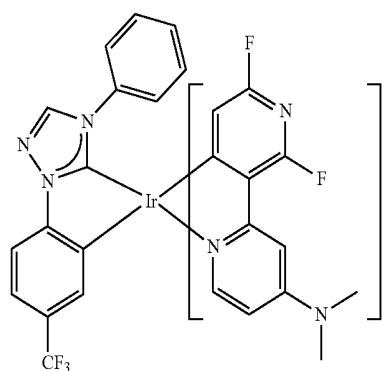
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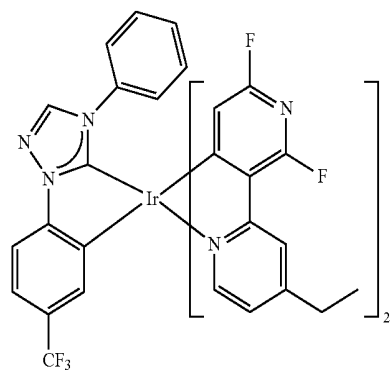
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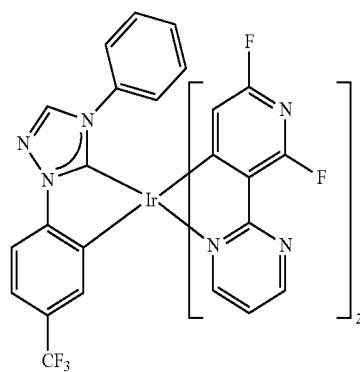
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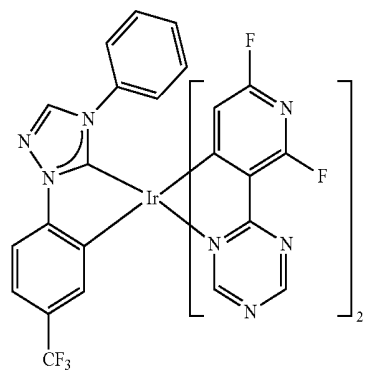


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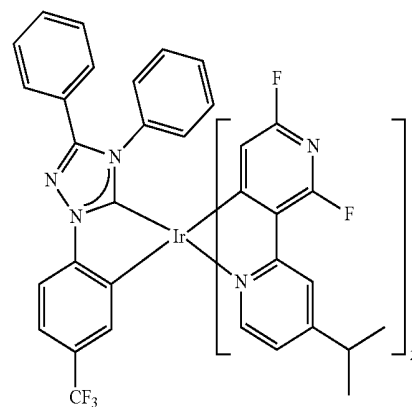
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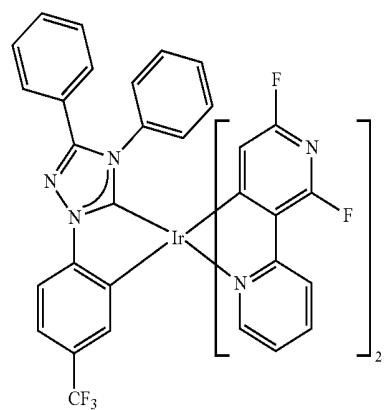
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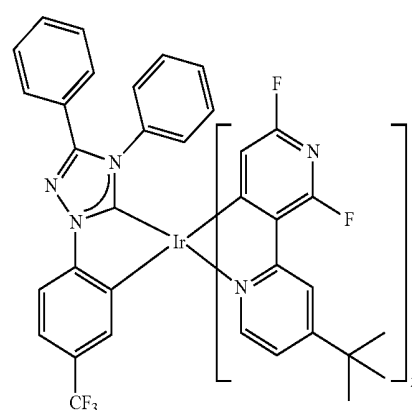
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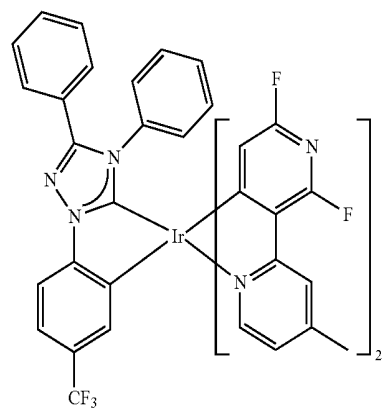
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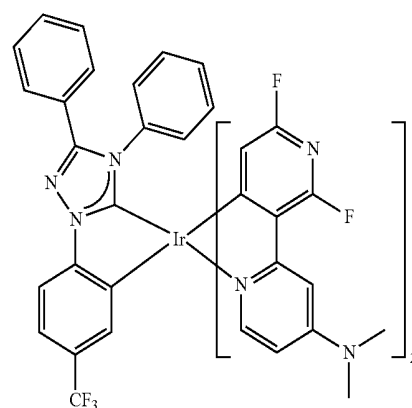


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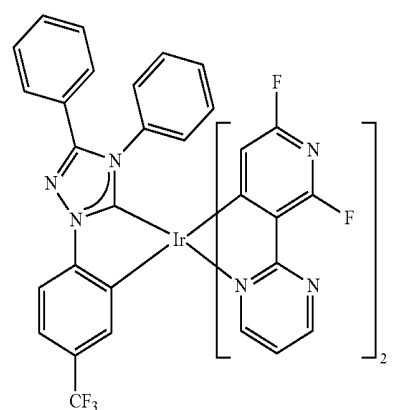
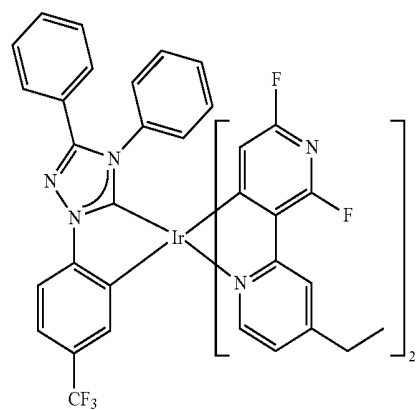
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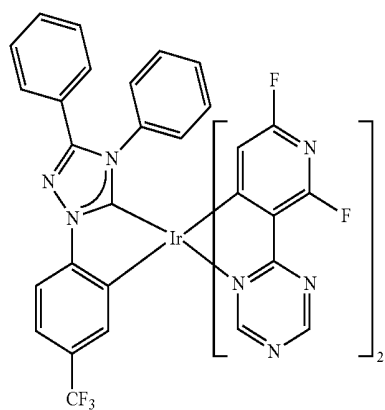
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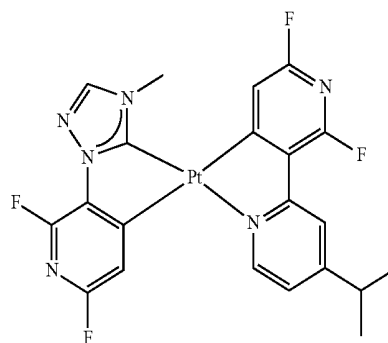


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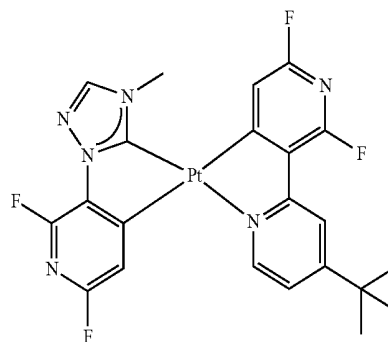
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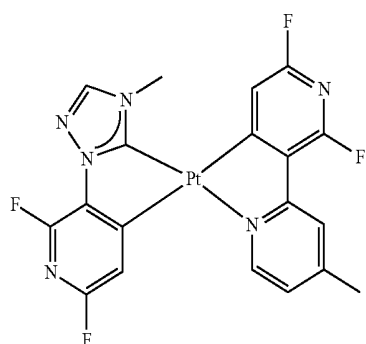
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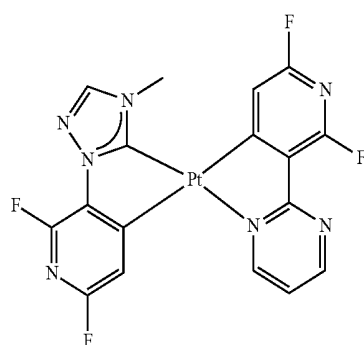
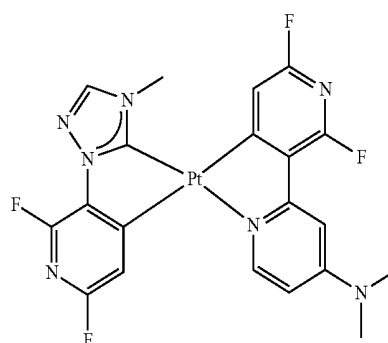
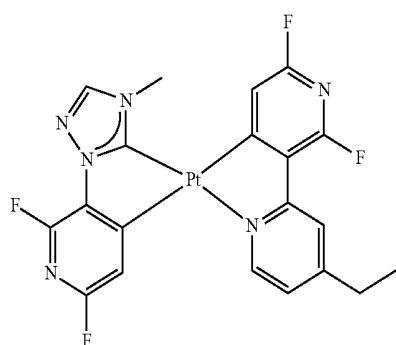
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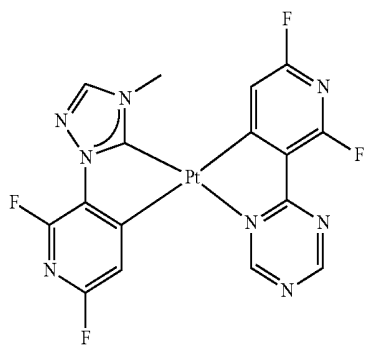


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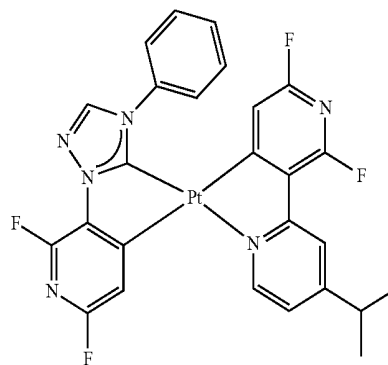


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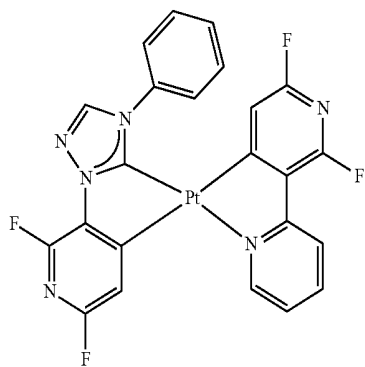


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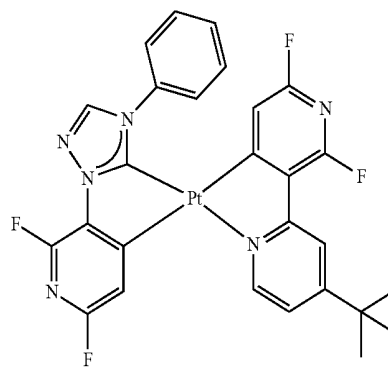
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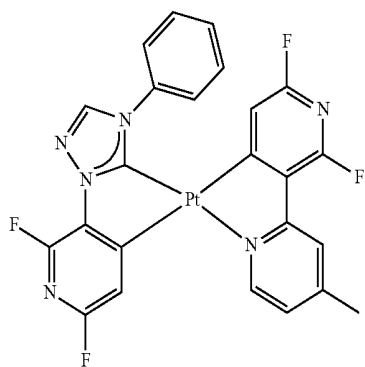
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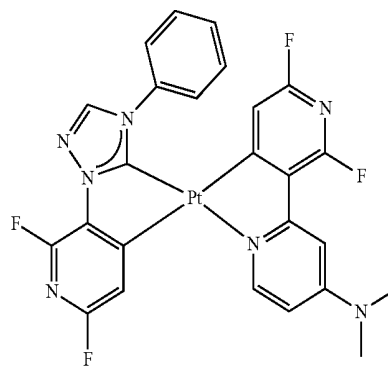
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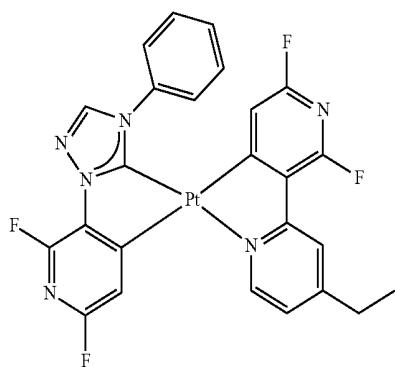
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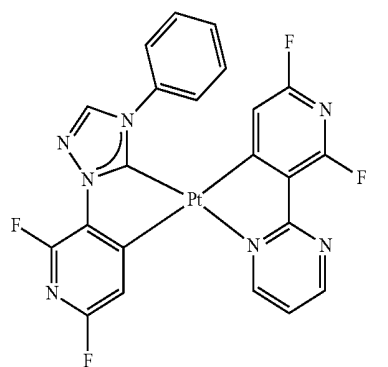
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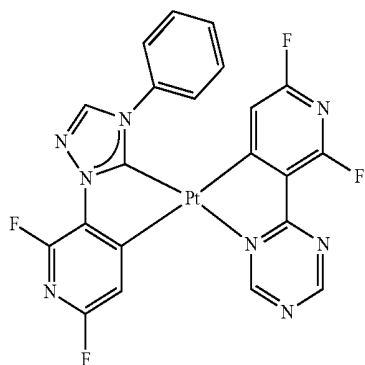


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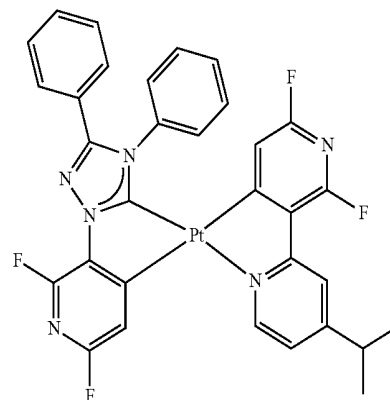


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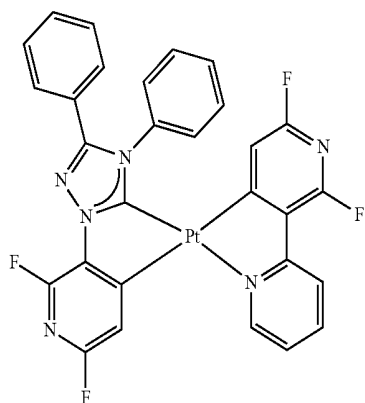
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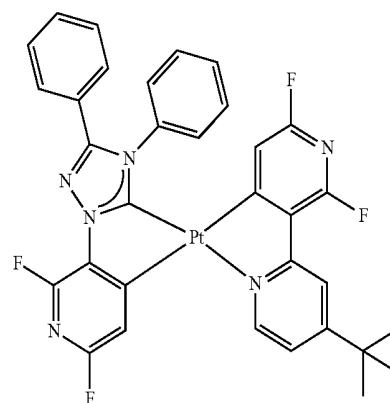
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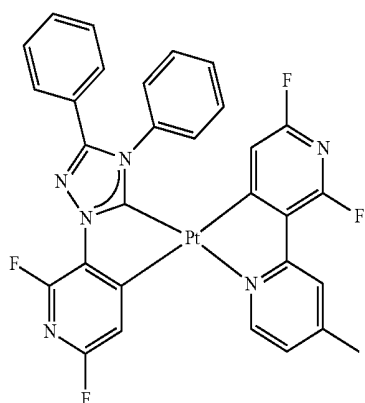
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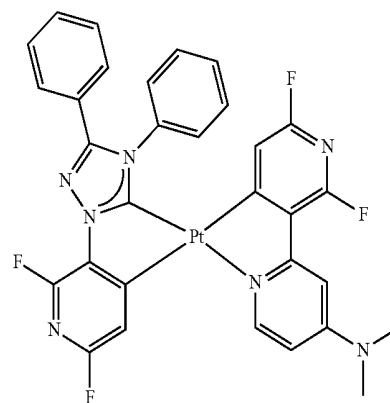
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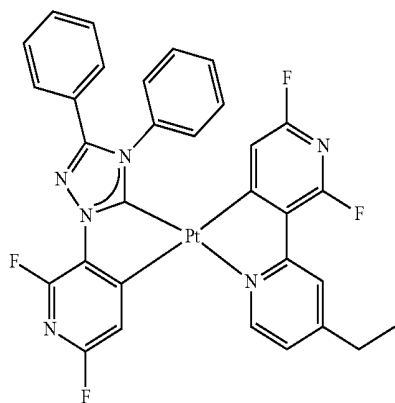
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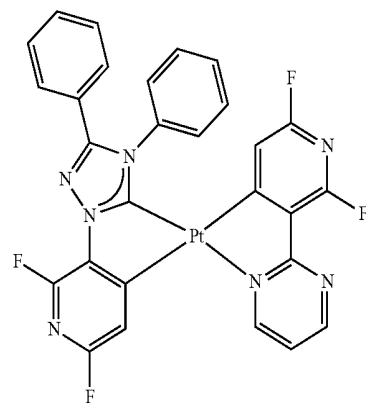
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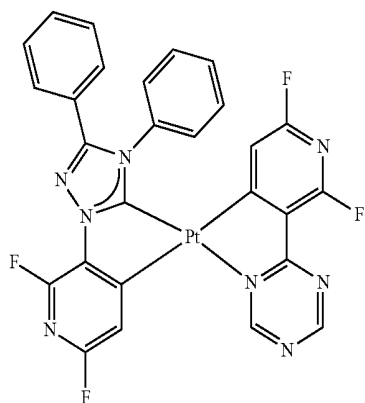


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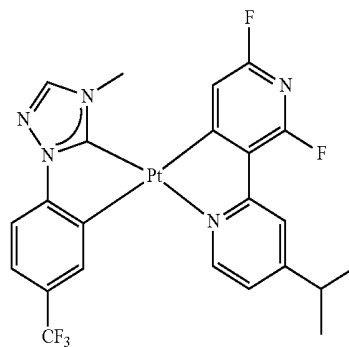
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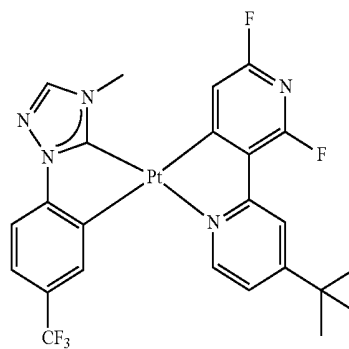
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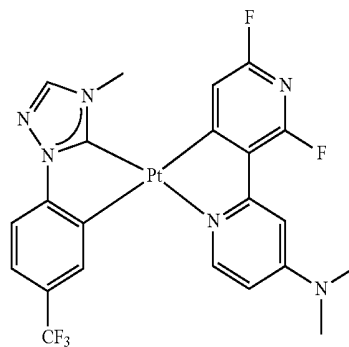
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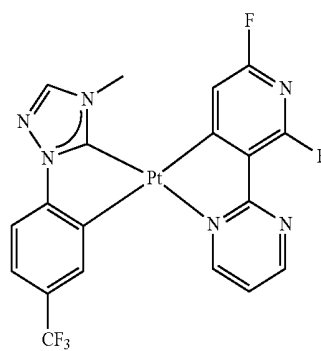
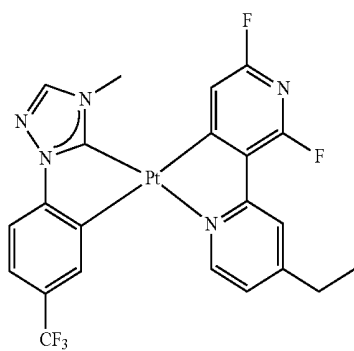
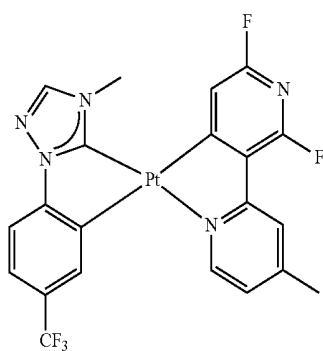
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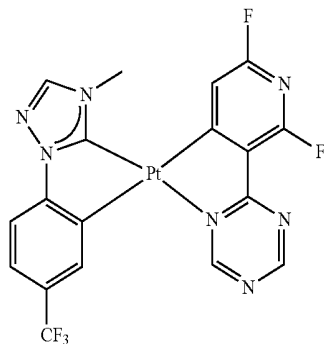


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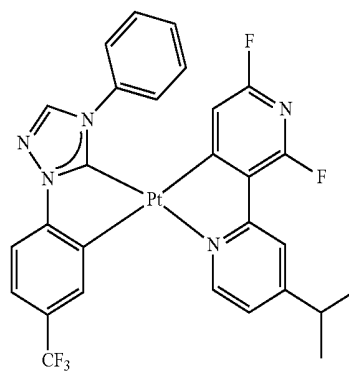


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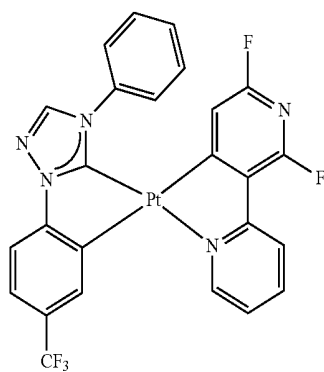


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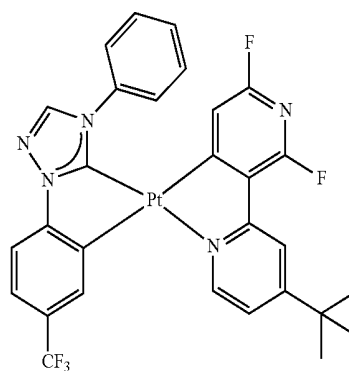
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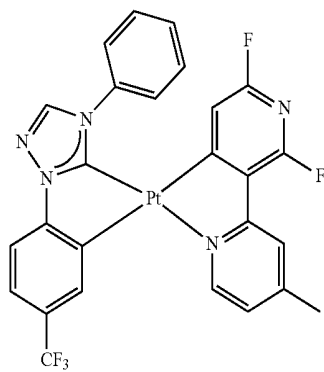
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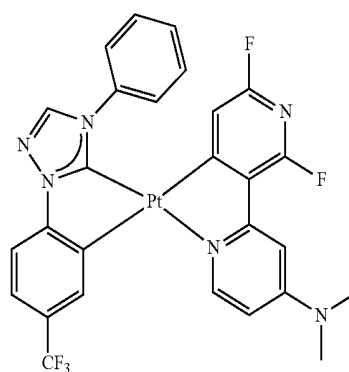
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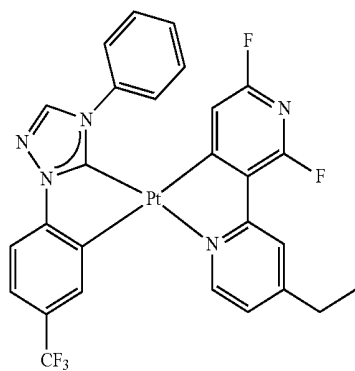
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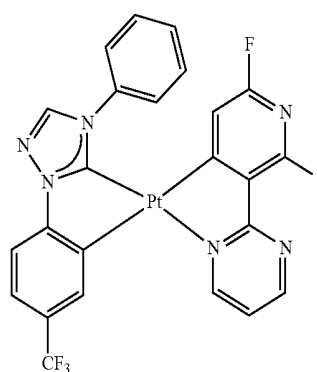
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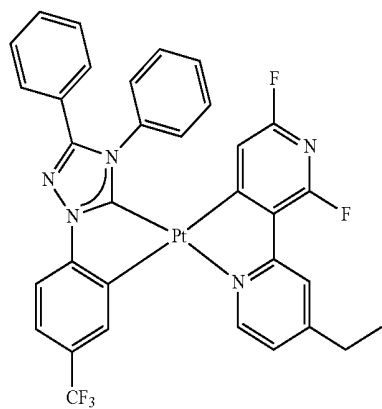
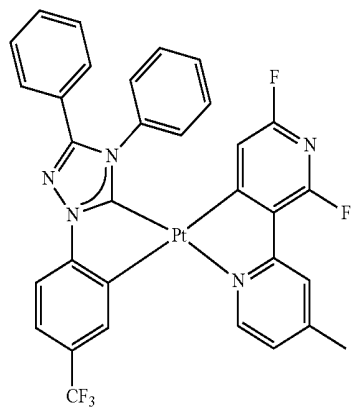
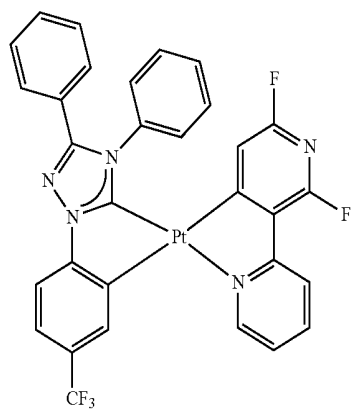
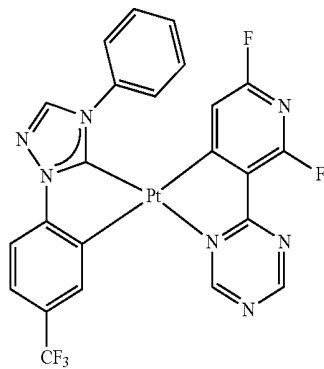


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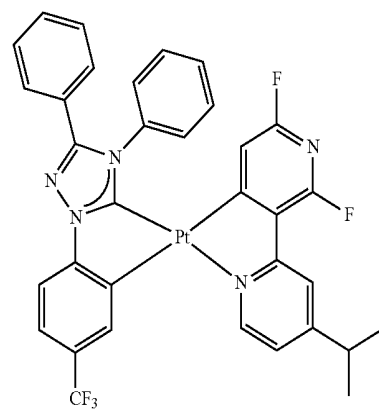
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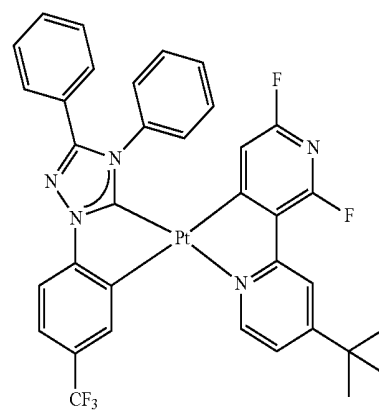
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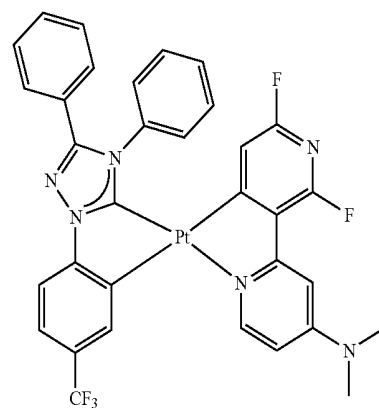
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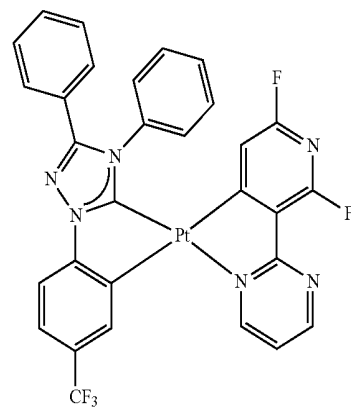
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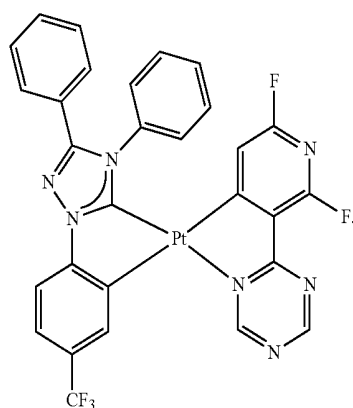
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95

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96



15. The organometallic compound of claim 1, wherein the organometallic compound emits blue light having a maximum emission wavelength of about 445 nm to about 465 nm.

16. 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, the organic layer including an emission layer, wherein the organic layer comprises at least one of the organometallic compound of claim 1.

17. The organic light-emitting device of claim 16, wherein the first electrode is an anode, the second electrode is a cathode, the organic layer further comprises a hole transport region between the first electrode and the emission layer and an electron transport region between the emission layer and the second electrode,

the hole transport region comprises a hole injection layer, a hole transport layer, an emission auxiliary layer, an electron blocking layer, or any combination thereof, and

the electron transport region comprises a hole blocking layer, an electron transport layer, an electron injection layer, or any combination thereof.

18. The organic light-emitting device of claim 16, wherein the organometallic compound is comprised in the emission layer, the emission layer further comprises a host, and an amount of the host in the emission layer is larger than an amount of the organometallic compound in the emission layer.

19. The organic light-emitting device of claim 18, wherein the host comprises at least one selected from a silicon-containing compound and a phosphine oxide-containing compound.

20. The organic light-emitting device of claim 17, wherein the hole transport region comprises a p-dopant, of which a lowest unoccupied molecular orbital (LUMO) energy level is lower than -3.5 eV.

* * * * *

专利名称(译)	有机金属化合物和包括其的有机发光器件		
公开(公告)号	US20180114927A1	公开(公告)日	2018-04-26
申请号	US15/789749	申请日	2017-10-20
[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	KO SOOBYUNG KIM SUNGBUM KIM YOUNGKOOK AHN HEECHOON JUN MIEUN HWANG SEOKHWAN		
发明人	KO, SOOBYUNG KIM, SUNGBUM KIM, YOUNGKOOK AHN, HEECHOON JUN, MIEUN HWANG, SEOKHWAN		
IPC分类号	H01L51/00		
CPC分类号	H01L51/0085 H01L51/0067 C07D213/61 C07D213/89 C07D249/10 H01L51/0087 H01L51/5016		
优先权	1020160136575 2016-10-20 KR		
外部链接	Espacenet USPTO		

摘要(译)

提供有机金属化合物和包含其的有机发光装置。有机金属化合物具有式1：其中M是第三行过渡金属，L1是由式2A表示的配体，n1是1,2或3，L2是由式表示的配体2B和n2为0,1或2.当n1为2或更多时，两个或更多个L1(s)彼此相同或不同。当n2为2或更多时，两个或更多个L2彼此相同或不同。提供了关于式2A和式2B的细节。

10

