

US010693076B2

(12) United States Patent
Hwang et al.**(10) Patent No.: US 10,693,076 B2****(45) Date of Patent: *Jun. 23, 2020****(54) CONDENSED-CYCLIC COMPOUND AND ORGANIC LIGHT-EMITTING DEVICE COMPRISING THE SAME****(71) Applicant: Samsung Display Co., Ltd.,** Yongin-si, Gyeonggi-do (KR)**(72) Inventors: Jinsoo Hwang,** Yongin-si (KR); **Myeongsuk Kim,** Yongin-si (KR); **Sungwook Kim,** Yongin-si (KR)**(73) Assignee: Samsung Display Co., Ltd.** (KR)**(*) Notice:** Subject to any disclaimer, the term of this patent is extended or adjusted under 35 U.S.C. 154(b) by 202 days.

This patent is subject to a terminal disclaimer.

(21) Appl. No.: 15/703,770**(22) Filed: Sep. 13, 2017****(65) Prior Publication Data**

US 2018/0076392 A1 Mar. 15, 2018

(30) Foreign Application Priority Data

Sep. 13, 2016 (KR) 10-2016-0118212

(51) Int. Cl.**H01L 51/00** (2006.01)**C07F 3/00** (2006.01)

(Continued)

(52) U.S. Cl.CPC **H01L 51/0058** (2013.01); **C07D 307/77**(2013.01); **C07D 307/91** (2013.01);

(Continued)

(58) Field of Classification Search

None

See application file for complete search history.

(56) References Cited

U.S. PATENT DOCUMENTS

8,901,298 B2 12/2014 Parham et al.
9,126,970 B2* 9/2015 Pflumm C07D 209/82
(Continued)

FOREIGN PATENT DOCUMENTS

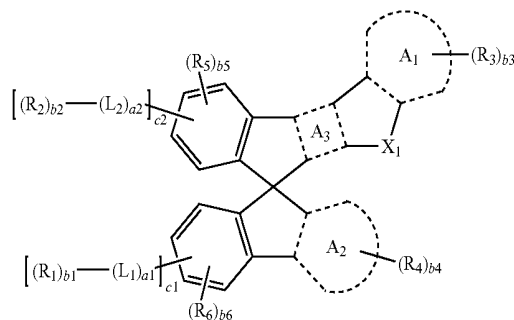
KR 10-2011-0088378 A 8/2011
KR 10-2013-0051807 A 5/2013

(Continued)

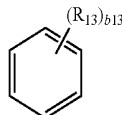
Primary Examiner — Robert S Loewe**(74) Attorney, Agent, or Firm** — Innovation Counsel LLP**(57) ABSTRACT**

A condensed-cyclic compound and an organic light-emitting device including the same are provided. The condensed-cyclic compound is represented by Formula 1:

Formula 1



Formula 2A



(Continued)

10**190****150****110**

-continued

Formula 2B



wherein, in Formulae 1, 2A, and 2B, ring A₁ is selected from a benzene group, a naphthalene group, a pyridine group, a pyrimidine group, a pyrazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, and a cinnoline group; ring A₂ is selected from a pyridine group, a pyrimidine group, a pyrazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, and a cinnoline group; and ring A₃ is selected from groups represented by Formula 2A and groups represented by Formula 2B. Further details about the compounds are presented.

20 Claims, 4 Drawing Sheets

(51) Int. Cl.

<i>C07D 307/77</i>	(2006.01)
<i>C07D 307/91</i>	(2006.01)
<i>C07D 403/04</i>	(2006.01)
<i>C07D 405/10</i>	(2006.01)
<i>C07D 405/12</i>	(2006.01)
<i>C07D 405/14</i>	(2006.01)
<i>C07D 409/04</i>	(2006.01)
<i>C07D 409/14</i>	(2006.01)
<i>C07D 471/10</i>	(2006.01)
<i>C07F 9/655</i>	(2006.01)
<i>C07F 9/6558</i>	(2006.01)
<i>H01L 51/50</i>	(2006.01)

(52) U.S. Cl.

CPC	<i>C07D 403/04</i> (2013.01); <i>C07D 405/10</i> (2013.01); <i>C07D 405/12</i> (2013.01); <i>C07D 405/14</i> (2013.01); <i>C07D 409/04</i> (2013.01); <i>C07D 409/14</i> (2013.01); <i>C07D 471/10</i> (2013.01); <i>C07F 3/006</i> (2013.01); <i>C07F</i>
-----------	--

9/65517 (2013.01); *C07F 9/65586* (2013.01); *H01L 51/0052* (2013.01); *H01L 51/0072* (2013.01); *H01L 51/0073* (2013.01); *H01L 51/5012* (2013.01); *H01L 51/5056* (2013.01); *H01L 51/5072* (2013.01); *H01L 51/5092* (2013.01); *H01L 51/5096* (2013.01); *H01L 51/0067* (2013.01); *H01L 2251/558* (2013.01)

(56)

References Cited

U.S. PATENT DOCUMENTS

9,288,869 B2	3/2016	Han et al.	
10,186,666 B2 *	1/2019	Kim	H01L 51/0067
2013/0240796 A1 *	9/2013	Parham	C07D 403/10 252/500
2014/0225040 A1	8/2014	Parham et al.	
2014/0225046 A1	8/2014	Jatsch et al.	
2015/0060808 A1 *	3/2015	Kim	H01L 51/0052 257/40
2015/0318478 A1	11/2015	Pflumm et al.	
2016/0043330 A1 *	2/2016	Kim	H01L 51/0085 257/40
2016/0126475 A1 *	5/2016	Lee	H01L 51/0072 257/40
2016/0126479 A1	5/2016	Hwang et al.	
2016/0190466 A1 *	6/2016	Pfister	H01L 51/0056 252/500
2016/0214942 A1 *	7/2016	Parham	C07D 401/14
2016/0218300 A1 *	7/2016	Xia	H01L 51/0072
2016/0293848 A1	10/2016	Kim et al.	
2016/0308147 A1 *	10/2016	Parham	C07F 15/0033

FOREIGN PATENT DOCUMENTS

KR	10-2013-0122602 A	11/2013	
KR	10-1381505 B1	3/2014	
KR	10-2014-0054132 A	5/2014	
KR	10-2014-0069199 A	6/2014	
KR	20150006802 A *	1/2015 C07D 487/10
KR	10-2015-0132660 A	11/2015	
KR	10-2016-0053423 A	5/2016	
KR	10-2016-0119904 A	10/2016	
WO	WO-2011025282 *	3/2011 C09K 11/06
WO	WO-2014104600 A1 *	7/2014 H05B 33/14
WO	WO-2016017919 A2 *	2/2016 C07D 307/91

* cited by examiner

FIG. 1

10

190
150
110

FIG. 2

20

190
150
110
210

FIG. 3

30

220
190
150
110

FIG. 4

40

220
190
150
110
210

1
CONDENSED-CYCLIC COMPOUND AND
ORGANIC LIGHT-EMITTING DEVICE
COMPRISING THE SAME

CROSS-REFERENCE TO RELATED
 APPLICATION

This application claims the benefit of Korean Patent Application No. 10-2016-0118212, filed on Sep. 13, 2016 in the Korean Intellectual Property Office, the disclosure of which is incorporated herein in its entirety by reference.

BACKGROUND

1. Field

One or more embodiments relate to a condensed-cyclic compound and an organic light-emitting device including the same.

2. Description of the Related Art

Organic light-emitting devices (OLEDs) are self-emission devices that have wide viewing angles, high contrast ratios, and short response times. In addition, the OLEDs exhibit high luminance, driving voltage, and response speed characteristics, and produce full-color images.

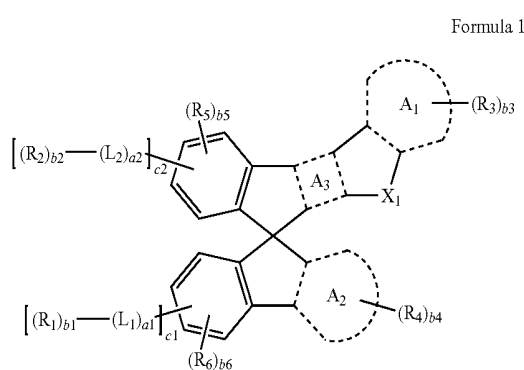
Organic light-emitting devices may include a first electrode disposed on a substrate. Organic light-emitting devices may include a hole transport region, an emission layer, an electron transport region, and a second electrode sequentially disposed on the first electrode. Holes provided from the first electrode may move toward the emission layer through the hole transport region. Electrons provided from the second electrode may move toward the emission layer through the electron transport region. Carriers, such as holes and electrons, may recombine in the emission layer to produce excitons. The excitons may transition from an excited state to a ground state, thus generating light.

SUMMARY

One or more embodiments include a condensed-cyclic compound and an organic light-emitting device including the same.

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

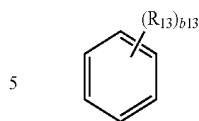
According to one or more embodiments, a condensed-cyclic compound is represented by Formula 1:



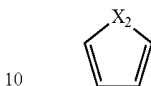
2

-continued

Formula 2A



Formula 2B



wherein, in Formulae 1, 2A, and 2B,

ring A_1 is selected from a naphthalene group, a pyridine group, a pyrimidine group, a pyrazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, and a cinnoline group,

ring A_2 is selected from a pyridine group, a pyrimidine group, a pyrazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, and a cinnoline group,

ring A_3 is selected from groups represented by Formula 2A and groups represented by Formula 2B,

X_1 is selected from N- $[(L_{11})_{a11}-(R_{11})_{b11}]$, O, and S,

X_2 is selected from N- $[(L_{12})_{a12}-(R_{12})_{b12}]$, O, and S,

L_1 and L_2 are each independently selected from substituted or unsubstituted condensed polycyclic groups, in which 3 or more carbocyclic groups are condensed with each other,

a_1 and a_2 are each independently an integer from 1 to 5; when a_1 is 2 or greater, at least two L_1 (s) are identical to or different from each other; when a_2 is 2 or greater, at least two L_2 (s) are identical to or different from each other,

L_{11} is selected from a substituted or unsubstituted C_3 - C_{10} cycloalkylene group, a substituted or unsubstituted C_3 - C_{10} cycloalkenylene group, a substituted or unsubstituted C_6 - C_{60} arylene group, and a substituted or unsubstituted divalent non-aromatic condensed polycyclic group,

a_{11} is an integer from 0 to 5,

L_{12} is 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,

a_{12} is an integer from 0 to 5,

R_{11} is selected from 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_6 - C_{60} aryl group, and a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group,

b_{11} is an integer from 1 to 4,

R_1 to R_6 , R_{12} , and R_{13} 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_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 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),

b1, b2, b5, b6, and b12 are each independently an integer from 0 to 4,

b3 and b4 are each independently an integer from 0 to 6, b13 is selected from 0, 1, and 2,

c1 and c2 are each independently an integer from 0 to 4, a sum of c1 and c2 is 1 or greater,

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

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, and a C₁-C₆₀ alkoxy group;

a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group, each substituted with at least one selected from 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);

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, a biphenyl group, and a terphenyl 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);

wherein 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 C₆-C₆₀ aryl group substituted with a C₁-C₆₀ alkyl group, a C₆-C₆₀ aryl group substituted with a C₆-C₆₀ aryl group, a terphenyl group, a C₁-C₆₀ heteroaryl group, a C₁-C₆₀ heteroaryl group substituted with a C₁-C₆₀ alkyl group, a C₁-C₆₀ heteroaryl group substituted with a C₆-C₆₀ aryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group.

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 disposed between the first electrode and the second electrode, wherein the organic layer includes an emission layer and at least one condensed-cyclic compound described above.

BRIEF DESCRIPTION OF THE DRAWINGS

These and/or other aspects will become apparent and more readily appreciated from the following description of the embodiments, taken in conjunction with the accompanying drawings in which:

FIG. 1 is a schematic cross-sectional view illustrating an organic light-emitting device according to an embodiment;

FIG. 2 is a schematic cross-sectional view illustrating an organic light-emitting device according to an embodiment;

FIG. 3 is a schematic cross-sectional view illustrating an organic light-emitting device according to an embodiment; and

FIG. 4 is a schematic cross-sectional view illustrating an organic light-emitting device according to an embodiment.

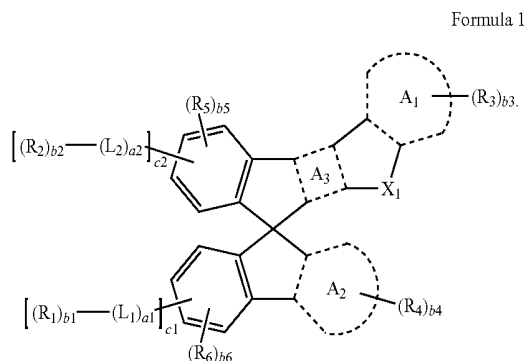
DETAILED DESCRIPTION

Reference will now be made in detail to embodiments, examples of which are illustrated in the accompanying drawings, wherein like reference numerals refer to like elements throughout. In this regard, the present embodiments may have different forms and should not be construed as being limited to the descriptions set forth herein. Accordingly, the embodiments are merely described below, by referring to the figures, to explain aspects of the present description. As used herein, the term “and/or” includes any and all combinations of one or more of the associated listed items. Expressions such as “at least one of,” when preceding

5

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

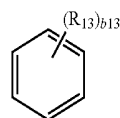
A condensed-cyclic compound may be represented by Formula 1:



Rings A_1 and A_2 in Formula 1 may each independently be fused to an adjacent 5-membered ring by sharing a carbon atom. In Formula 1, ring A_1 may be selected from a benzene group, a naphthalene group, a pyridine group, a pyrimidine group, a pyrazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, and a cinnoline group, and ring A_2 may be selected from a pyridine group, a pyrimidine group, a pyrazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, and a cinnoline group.

In some embodiments, in Formula 1, ring A_1 may be selected from a benzene group and a naphthalene group, and ring A_2 may be selected from a pyridine group, a quinoline group, and an isoquinoline group.

In Formula 1, ring A_3 may be fused to two adjacent 5-membered rings by sharing carbon atoms. In Formula 1, ring A_3 may be selected from groups represented by Formula 2A and groups represented by Formula 2B.



R_{13} and b_{13} in Formula 2A and X_2 in Formula 2B may be the same as those described herein.

In one embodiment, in Formula 1, ring A_3 may be a group represented by Formula 2A, but embodiments are not limited thereto.

X_1 in Formula 1 may be $N-[(L_{11})_{a11}-(R_{11})_{b11}]$, O, or S, and X_2 in Formula 2B may be $N-[(L_{12})_{a12}-(R_{12})_{b12}]$, O, or S.

In some embodiments, in Formula 1, X_1 may be $N-[(L_{11})_{a1}-(R_{11})_{b11}]$.

In Formula 1, L_1 and L_2 may each independently be selected from substituted or unsubstituted condensed polycyclic groups, in which 3 or more carbocyclic groups may be condensed with each other. L_1 and L_2 may include carbon atoms as ring-forming atoms, but may not include heteroatoms (e.g., N, O, S, or P) as ring forming-atoms. Therefore,

6

for example, a naphthylene group may not be L_1 or L_2 because a naphthylene group is a condensed polycyclic group in which two carbocyclic groups are condensed; and a pyridinylene group may not be L_1 or L_2 because a pyridinylene group contains N as a ring-forming atom.

In one embodiment, in Formula 1, L_1 and L_2 may each independently be selected from

an indacenylene group, an acenaphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzofluorenylene group, a dibenzofluorenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylene group, a fluoranthenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a naphthacenylene group, a picenylene group, a perylenylene group, a pentaphenylene group, a hexacenylene group, a pentacenylene group, a rubicenylene group, a coronenylene group, and an ovalenylene group; and

an indacenylene group, an acenaphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzofluorenylene group, a dibenzofluorenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylene group, a fluoranthenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a naphthacenylene group, a picenylene group, a perylenylene group, a pentaphenylene group, a hexacenylene group, a pentacenylene group, a rubicenylene group, a coronenylene group, and an ovalenylene 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 cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, 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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a 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 dibenzothiofenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a thiadiazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and $-Si(Q_{31})(Q_{32})(Q_{33})$,

wherein 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.

In one embodiment, in Formula 1, L_1 and L_2 may each independently be selected from

a phenalenylylene group, a phenanthrenylylene group, an anthracenylylene group, a fluoranthenylylene group, a triphenylylylene group, a pyrenylylene group, a chrysenylylene group, and a perylynylylene group; and

a phenalenylylene group, a phenanthrenylylene group, an anthracenylylene group, a fluoranthenylylene group, a triphenylylylene group, a pyrenylylene group, a chrysenylylene group, and a perylynylylene 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 cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, 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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenylylene group, a phenanthrenylylene group, an anthracenylylene group, a fluoranthenylylene group, a triphenylylylene group, a pyrenyl group, a chrysenyl group, a naphthacenylylene group, a picenyl group, a perylynylylene group, a pentaphenyl group, a hexacenylylene group, a pentacenylylene group, a rubicenylylene group, a coronenylylene group, a pyrolylylene group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a benzimidazolyl group, a benzofuran 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 dibenzofuran group, a dibenzothiophenyl group, a benzocarbazolyl group, a dibenzocarbazolyl group, a thiadiazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and —Si(Q_{31})(Q_{32})(Q_{33}),

wherein 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.

In Formula 1, a_1 and a_2 may each independently be an integer from 1 to 5; when a_1 is 2 or greater, at least two L_1 (s) may be identical to or different from each other; and when a_2 is 2 or greater, at least two L_2 (s) may be identical to or different from each other. That is, in Formula 1, L_1 may be essentially present in a group represented by $*[(L_1)_{a_1}-(R_1)_{b_1}]$, and L_2 may be essentially present in a group represented by $*[(L_2)_{a_2}-(R_2)_{b_2}]$.

In one embodiment, in Formula 1, a_1 and a_2 may each independently be 1 or 2. In some embodiments, in Formula 1, a_1 and a_2 may each independently be 1, but embodiments are not limited thereto.

In Formulae 1, 2A, and 2B, L_{11} may be selected from a substituted or unsubstituted C_3 - C_{10} cycloalkylene group, a substituted or unsubstituted C_3 - C_{10} cycloalkenylene group, a substituted or unsubstituted C_6 - C_{60} arylene group, and a substituted or unsubstituted divalent non-aromatic condensed polycyclic group, and

L_{12} 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.

In some embodiments, L_{11} may be selected from a phenylene group, a pentalenylylene group, an indenylene group, a naphthylene group, an azulenylylene group, a heptalenylene group, an indacenylene group, an acenaphthylene group, a fluorenylylene group, a spiro-fluorenylylene group, a benzofluorenylylene group, a dibenzofluorenylylene group, a phenalenylylene group, a phenanthrenylene group, an anthracenylylene group, a fluoranthenylylene group, a triphenylylylene group, a pyrenylene group, a chrysenylene group, a naphthacenylylene group, a picenylene group, a perylynylylene group, a pentaphenylene group, a hexacenylylene group, a pentacenylylene group, a rubicenylylene group, a coronenylylene group, and an ovalenylylene group; and

a phenylene group, a pentalenylylene group, an indenylene group, a naphthylene group, an azulenylylene group, a heptalenylene group, an indacenylene group, an acenaphthylene group, a fluorenylylene group, a spiro-fluorenylylene group, a benzofluorenylylene group, a dibenzofluorenylylene group, a phenalenylylene group, a phenanthrenylene group, an anthracenylylene group, a fluoranthenylylene group, a triphenylylylene group, a pyrenylene group, a chrysenylene group, a naphthacenylylene group, a picenylene group, a perylynylylene group, a pentaphenylene group, a hexacenylylene group, a pentacenylylene group, a rubicenylylene group, a coronenylylene group, and an ovalenylylene 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 cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, 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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenylylene group, a phenanthrenylene group, an anthracenylylene group, a fluoranthenylylene group, a triphenylylylene group, a pyrenyl group, a chrysenyl group, a naphthacenylylene group, a picenyl group, a perylynylylene group, a pentaphenyl group, a hexacenylylene group, a pentacenylylene group, a rubicenylylene group, a coronenylylene group, an ovalenylylene group, and —Si(Q_{31})(Q_{32})(Q_{33}),

wherein 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, and

L_{12} may be selected from a phenylene group, a pentalenylylene group, an indenylene group, a naphthylene group, an azulenylylene group, a heptalenylene group, an indacenylene group, an acenaphthylene group, a fluorenylylene group, a spiro-fluorenylylene group, a benzofluorenylylene group, a dibenzofluorenylylene group, a phenalenylylene group, a phenanthrenylene group, an anthracenylylene group, a fluoranthenylylene group, a triphenylylylene group, a pyrenylene group, a chrysenylene group, a naphthacenylylene group, a picenylene group, a perylynylylene group, a pentaphenylene group, a hexacenylylene group, a pentacenylylene group, a

rubicenylene group, a coronenylene group, an ovalenylene group, a pyrrolylene group, a thiophenylene group, a furanylene group, an imidazolylene group, a pyrazolylene group, a thiazolylene group, an isothiazolylene group, an oxazolylene group, an isoxazolylene group, a pyridinylene group, a pyrazinylene group, a pyrimidinylene group, a pyridazinylene group, an isoindolylene group, an indolylene group, an indazolylene group, a purinylene 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 carbazolylene group, a phenanthridinylene group, an acridinylene group, a phenanthrolinylene group, a phenazinylene group, a benzimidazolylene group, a benzofuranylene group, a benzothiophenylene group, an isobenzothiazolylene group, a benzoxazolylene group, an isobenzoxazolylene group, a triazolylene group, a tetrazolylene group, an oxadiazolylene group, a triazinylene group, a dibenzofuranylene group, a dibenzothiophenylene group, a benzocarbazolylene group, a dibenzocarbazolylene group, a thiadiazolylene group, an imidazopyridinylene group, and an imidazopyrimidinylene group; and

a phenylene group, a pentalenylene group, an indenylene group, a naphthylene group, an azulenylene group, a heptalenylene group, an indacenylene group, an acenaphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzofluorenylene group, a dibenzofluorenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylene group, a fluoranthenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a naphthacenylenylene group, a picenylene group, a perylenylene group, a pentaphenylenylene group, a hexacenylenylene group, a pentacenylenylene group, a rubicenylene group, a coronenylene group, an ovalenylene group, a pyrrolylene group, a thiophenylene group, a furanylene group, an imidazolylene group, a pyrazolylene group, a thiazolylene group, an isothiazolylene group, an oxazolylene group, an isoxazolylene group, a pyridinylene group, a pyrazinylene group, a pyrimidinylene group, a pyridazinylene group, an isoindolylene group, an indolylene group, an indazolylene group, a purinylene 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 carbazolylene group, a phenanthridinylene group, an acridinylene group, a phenanthrolinylene group, a phenazinylene group, a benzimidazolylene group, a benzofuranylene group, a benzothiophenylene group, an isobenzothiazolylene group, a benzoxazolylene group, an isobenzoxazolylene group, a triazolylene group, a tetrazolylene group, an oxadiazolylene group, a triazinylene group, a dibenzofuranylene group, a dibenzothiophenylene group, a benzocarbazolylene group, a dibenzocarbazolylene group, a thiadiazolylene group, an imidazopyridinylene group, and an imidazopyrimidinylene group, each substituted with at least one selected from 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 cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, 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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyre-

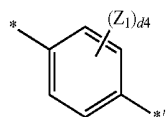
nyl group, a chrysenyl group, a naphthaceny group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexaceny group, a pentaceny group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a 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, a thiadiazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and —Si(Q₃₁)(Q₃₂)(Q₃₃),

wherein 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.

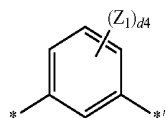
In one embodiment, L₁ and L₂ may each independently be selected from groups represented by Formulae 3-8, 3-9, 3-25, and 3-35 to 3-41,

L₁₁ may be selected from groups represented by Formulae 3-1 to 3-9, 3-25, and 3-33 to 3-41, wherein Y₁ in Formulae 3-3 and 3-4 may be C(Z₃)(Z₄), and

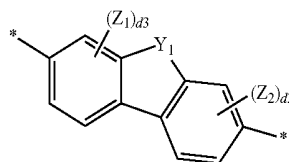
L₁₂ may be selected from groups represented by Formulae 3-1 to 3-41:



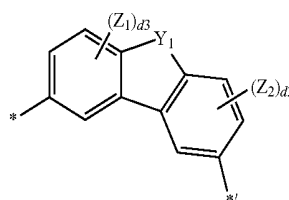
Formula 3-1



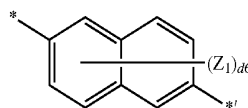
Formula 3-2



Formula 3-3

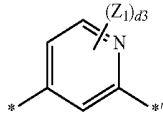
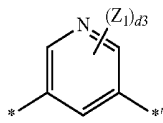
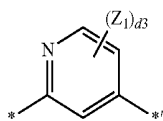
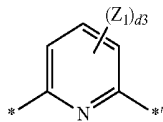
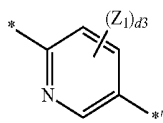
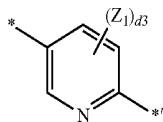
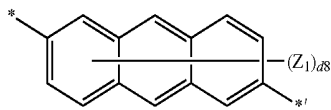
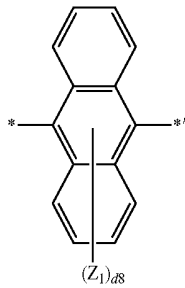
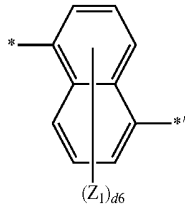
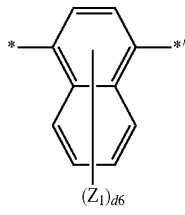


Formula 3-4



Formula 3-5

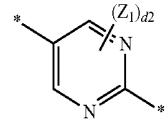
11
-continued



12
-continued

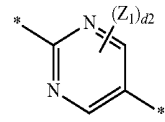
Formula 3-6

5

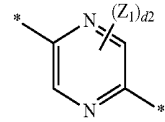


Formula 3-7

10

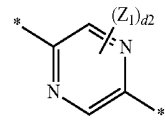


15

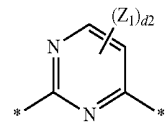


Formula 3-8

20

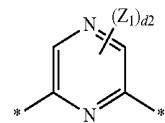


25



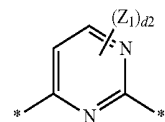
Formula 3-9

30



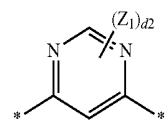
Formula 3-10

35



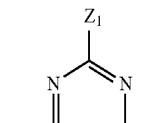
Formula 3-11

40



45

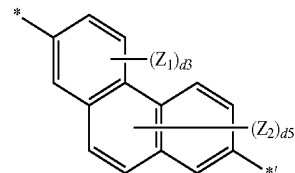
Formula 3-12



50

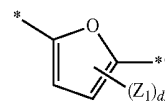
Formula 3-13

55



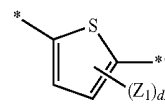
Formula 3-14

60



Formula 3-15

65



Formula 3-16

Formula 3-17

Formula 3-18

Formula 3-19

Formula 3-20

Formula 3-21

Formula 3-22

Formula 3-23

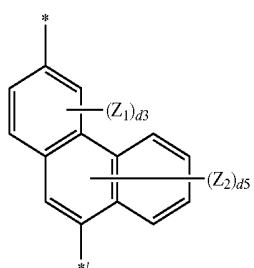
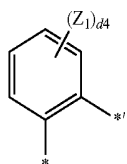
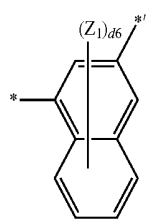
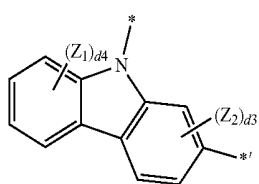
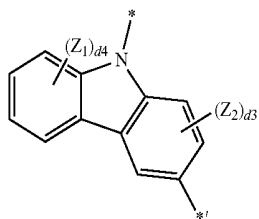
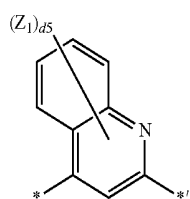
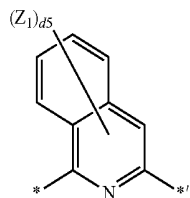
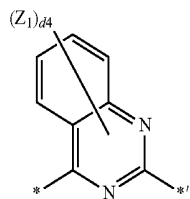
Formula 3-24

Formula 3-25

Formula 3-26

Formula 3-27

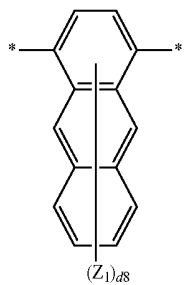
13
-continued



14
-continued

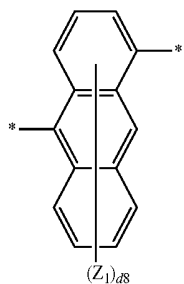
Formula 3-28

5



Formula 3-29 10

15

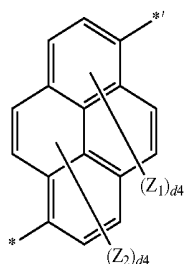


Formula 3-30 20

25

Formula 3-31

30

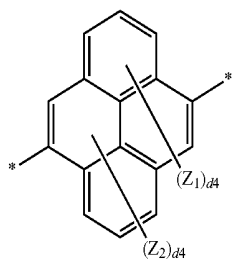


Formula 3-32 35

40

Formula 3-33

45



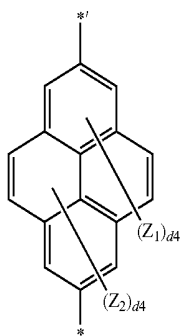
Formula 3-34 50

55

Formula 3-35

60

65



Formula 3-36

Formula 3-37

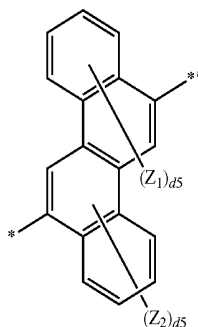
Formula 3-38

Formula 3-39

Formula 3-40

15

-continued



wherein, in Formulae 3-1 to 3-41,

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

Z_1 to Z_7 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 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 pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a dibenzofuranyl group, a dibenzothiophenyl group, a triazinyl group, a benzimidazolyl group, a phenanthrolinyl group, and $-Si(Q_{31})(Q_{32})(Q_{33})$,

wherein 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,

d_2 may be an integer selected from 1 and 2,

d_3 may be an integer from 1 to 3,

d_4 may be an integer from 1 to 4,

d_5 may be an integer from 1 to 5,

d_6 may be an integer from 1 to 6,

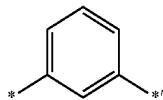
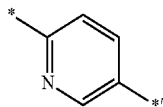
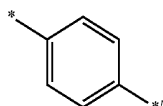
d_8 may be an integer from 1 to 8, and

* and *' each indicate a binding site to an adjacent atom.

In some embodiments, L_1 and L_2 may each independently be selected from groups represented by Formulae 4-11, 4-13, 4-27, and 4-29 to 4-35,

L_{11} may be selected from groups represented by Formulae 4-1, 4-3, 4-5, 4-7 to 4-13, 4-17, and 4-24 to 4-35, and

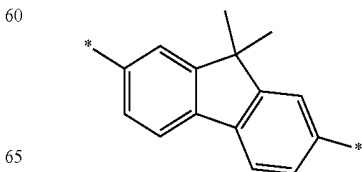
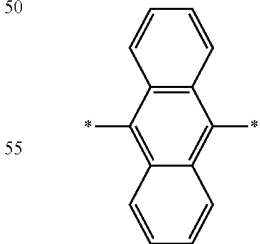
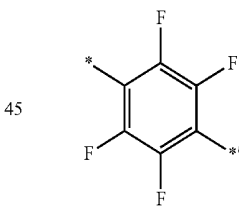
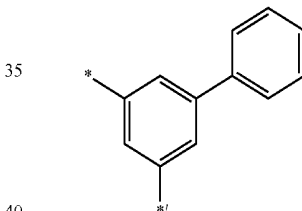
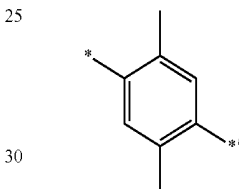
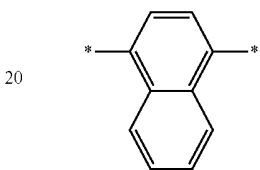
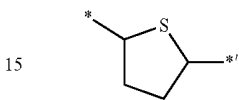
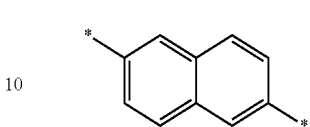
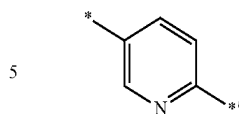
L_{12} may be selected from groups represented by Formulae 4-1 to 4-35, but embodiments are not limited thereto:



Formula 3-41

16

-continued



Formula 4-4

Formula 4-5

Formula 4-6

Formula 4-7

Formula 4-8

Formula 4-9

Formula 4-10

Formula 4-11

Formula 4-12

Formula 4-1

Formula 4-2

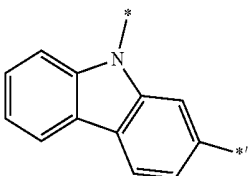
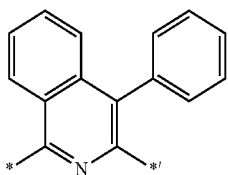
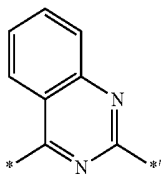
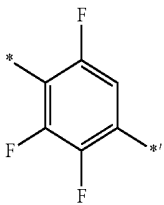
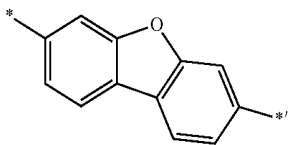
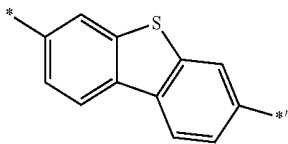
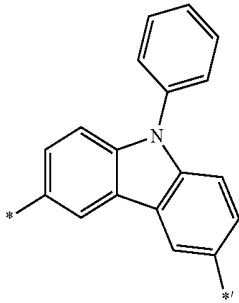
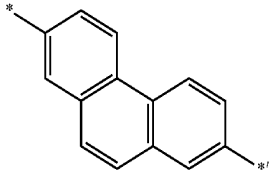
Formula 4-3

55

60

65

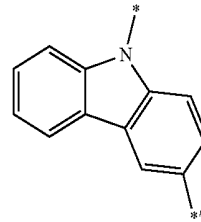
17
-continued



18
-continued

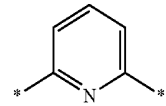
Formula 4-13

5

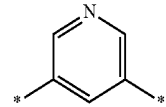


Formula 4-14 10

15

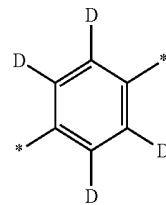


20



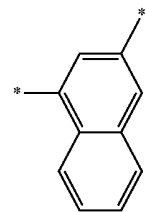
Formula 4-15

25



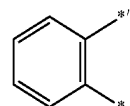
Formula 4-16

30



Formula 4-17

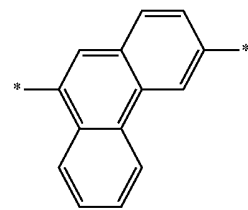
35



40

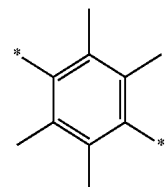
Formula 4-18

45



Formula 4-19

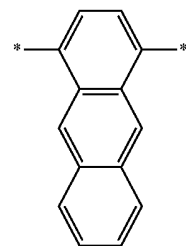
50



55

Formula 4-20

60



65

Formula 4-21

Formula 4-22

Formula 4-23

Formula 4-24

Formula 4-25

Formula 4-26

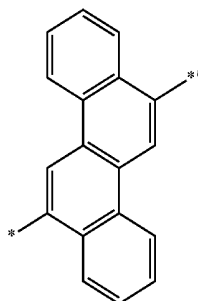
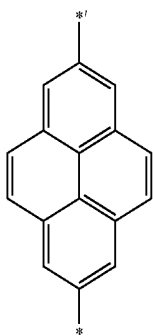
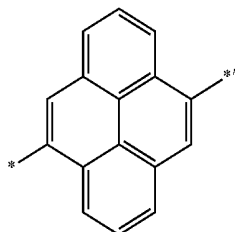
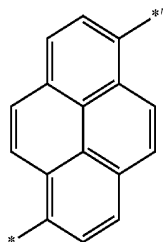
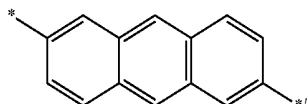
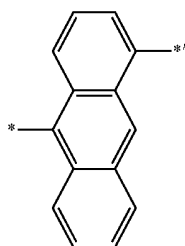
Formula 4-27

Formula 4-28

Formula 4-29

19

-continued



wherein, in Formulae 4-1 to 4-35, * and *' each indicate a binding site to an adjacent atom.

In the above formulae, a₁₁ and a₁₂ may each independently be an integer from 0 to 5; when a₁₁ is 2 or greater, at least two L₁₁ (s) may be identical to or different from each other; and when a₁₂ is 2 or greater, at least two L₂ (s) may be identical to or different from each other.

20

Formula 4-30

In one embodiment, in the above formulae, a₁₁ and a₁₂ may each independently be 0, 1, or 2. In some embodiments, in the above formulae, a₁₁ and a₁₂ may each independently be 0 or 1, but embodiments are not limited thereto.

5 In the above formulae, R₁₁ may be selected from a substituted or unsubstituted C₃-C₁₀ cycloalkyl group, a substituted or unsubstituted C₃-C₁₀ cycloalkenyl group, a substituted or unsubstituted C₆-C₆₀ aryl group, and a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group.

Formula 4-31

10 In some embodiments, R₁₁ may 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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthrenyl 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, and an ovalenyl group;

Formula 4-32

Formula 4-33

Formula 4-34

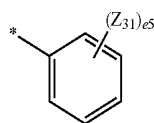
Formula 4-35

25 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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthrenyl 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, and an ovalenyl 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 cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, 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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthrenyl 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, and —Si(Q₃₁)(Q₃₂)(Q₃₃),

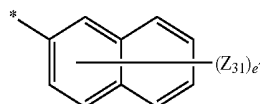
30 wherein Q₃₁ to Q₃₃ may each independently be selected from a C₁-C₁₀ alkyl group, a C₁-C₁₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group.

35 In some embodiments, R₁₁ may be selected from groups represented by Formulae 5-1 to 5-19:

Formula 5-1



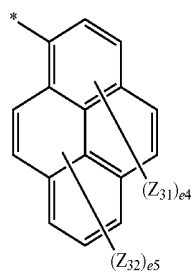
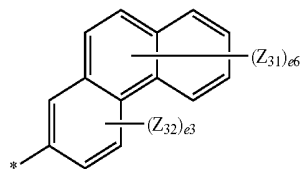
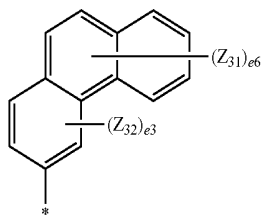
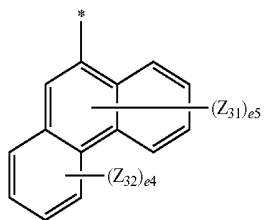
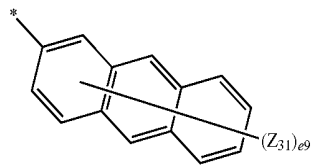
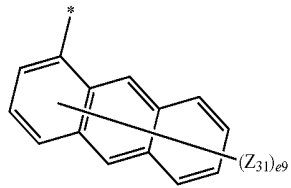
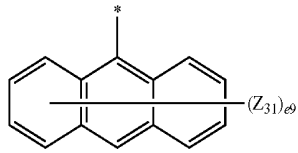
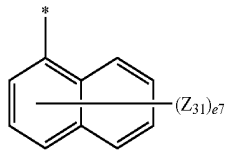
Formula 5-2



60

65

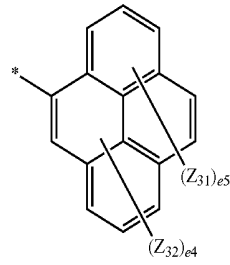
21
-continued



22
-continued

Formula 5-3

5

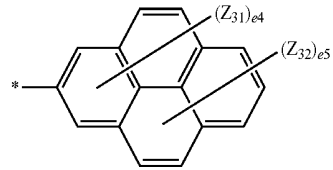


Formula 5-4

10

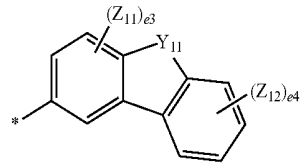
Formula 5-5

15



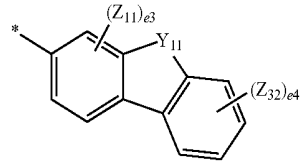
Formula 5-6

25



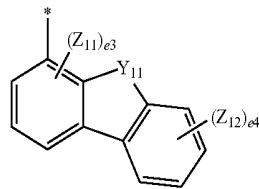
Formula 5-7

30



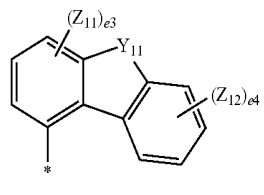
Formula 5-8

40



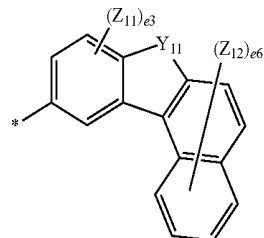
Formula 5-9

50

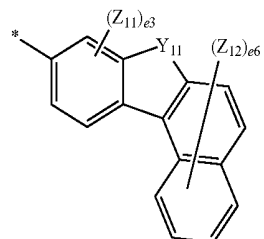


Formula 5-10

55



65



Formula 5-11

Formula 5-12

Formula 5-13

Formula 5-14

Formula 5-15

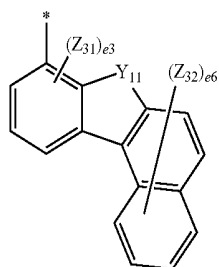
Formula 5-16

Formula 5-17

Formula 5-18

23

-continued



wherein, in Formulae 5-1 to 5-19,

Y_{11} may be $C(Z_{13})(Z_{14})$,

Z_{11} to Z_{14} 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 phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a biphenyl group, a terphenyl group, and $-Si(Q_{31})(Q_{32})(Q_{33})$,

wherein 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,

e_3 may be an integer from 1 to 3,

e_4 may be an integer from 1 to 4,

e_5 may be an integer from 1 to 5,

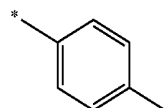
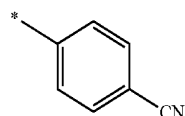
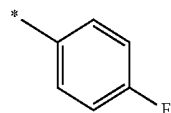
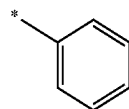
e_6 may be an integer from 1 to 6,

e_7 may be an integer from 1 to 7,

e_9 may be an integer from 1 to 9, and

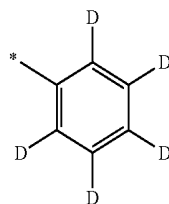
* indicates a binding site to an adjacent atom.

In some embodiments, R_{11} may be selected from groups represented by Formulae 6-1 to 6-27:

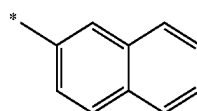


Formula 5-19

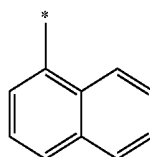
5



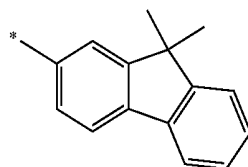
10



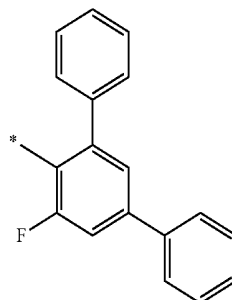
15



20



25



30

35

40

Formula 6-5

Formula 6-6

Formula 6-7

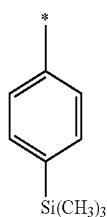
Formula 6-8

Formula 6-9

Formula 6-10

Formula 6-1

45



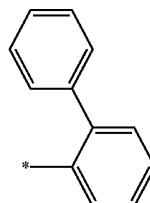
50

Formula 6-2

55

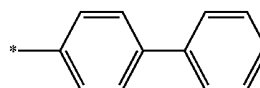
Formula 6-3

60



Formula 6-4

65

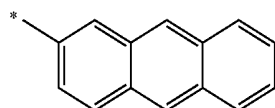
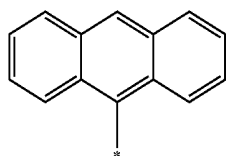
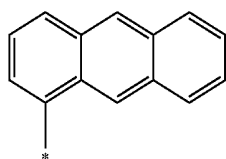
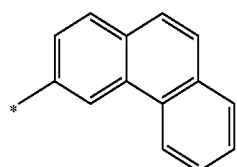
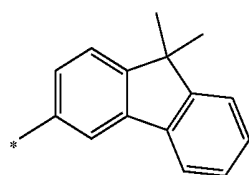
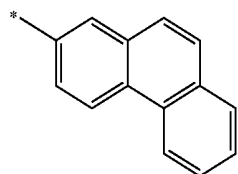
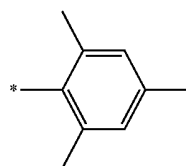
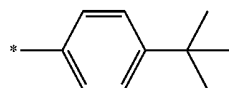
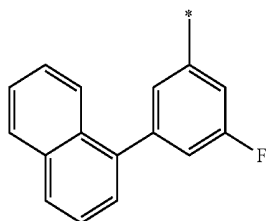


Formula 6-11

Formula 6-12

25

-continued

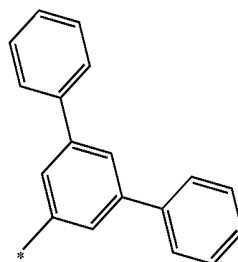


26

-continued

Formula 6-13

5

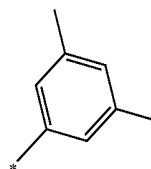


Formula 6-14

10

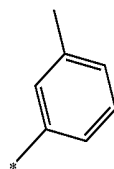
Formula 6-15

15



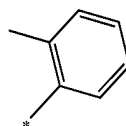
Formula 6-16

20



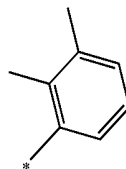
Formula 6-17

25



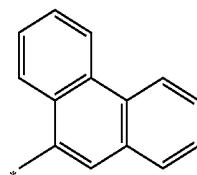
Formula 6-18

30



Formula 6-19

35



Formula 6-20

40

Formula 6-21

45

Formula 6-22

Formula 6-23

Formula 6-24

Formula 6-25

Formula 6-26

Formula 6-27

wherein, in Formulae 6-1 to 6-27, * indicates a binding site to an adjacent atom.

In the above formulae, b11 may be an integer from 1 to 4. In some embodiments, in the above formulae, b11 may be 1 or 2.

In some embodiments, in the above formulae, b11 may be 1, but embodiments are not limited thereto.

In the above formulae, R₁ to R₆, R₁₂, and R₁₃ 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 substi-

tuted 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).

In some embodiments, in the above formulae, R₁ to R₆, R₁₂, and R₁₃ 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, and a C₁-C₂₀ alkoxy group;

a C₁-C₂₀ alkyl group and a C₁-C₂₀ 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 phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group;

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

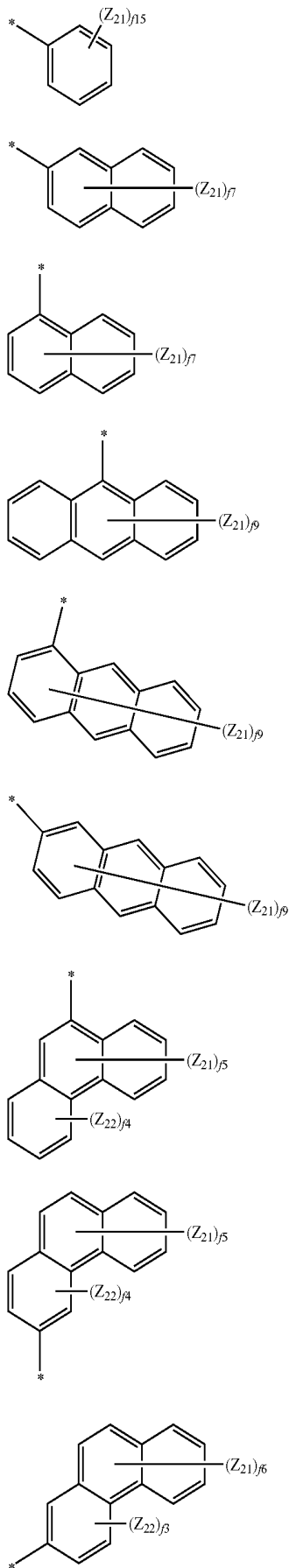
a phenyl group, a biphenyl group, a terphenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenyl group, a heptalenyl group, an indacenyl group, an acenaphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a

benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a 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, a dibenzosilolyl group, a thiadiazolyl group, an imidazopyridinyl group, and an imidazopyrimidinyl 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 cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, 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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a 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, a thiadiazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and —Si(Q₃₁)(Q₃₂)(Q₃₃); and

—Si(Q₁)(Q₂)(Q₃),

wherein Q₁ to Q₃ and 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.

In one embodiment, in the above formulae, R₁ to R₆, R₁₂, and R₁₃ 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-C₁₀ alkyl group, a C₁-C₁₀ alkoxy group, groups represented by Formulae 7-1 to 7-75, and —Si(Q₁)(Q₂)(Q₃), wherein Q₁ to Q₃ 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:



Formula 7-1

5

Formula 7-2

10

Formula 7-3

15

Formula 7-4

20

Formula 7-5

30

Formula 7-6

35

Formula 7-7

45

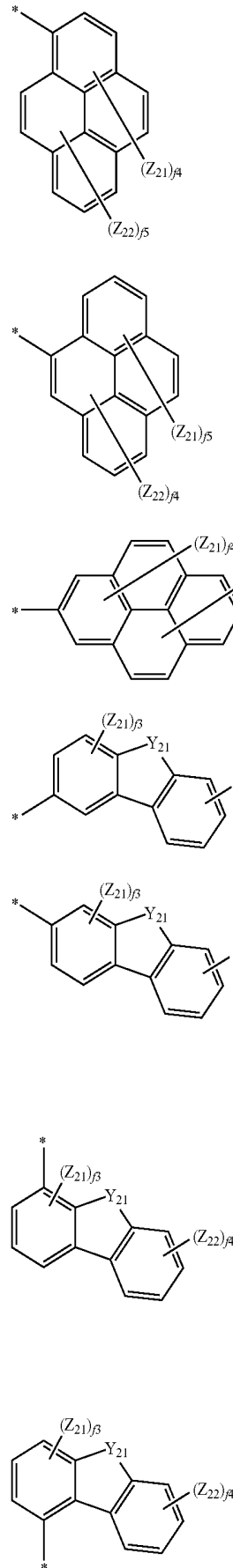
Formula 7-8

55

Formula 7-9

60

65



Formula 7-10

Formula 7-11

Formula 7-12

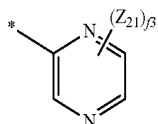
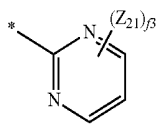
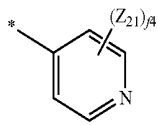
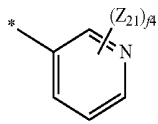
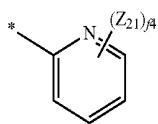
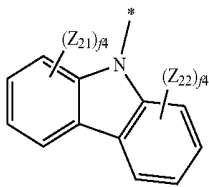
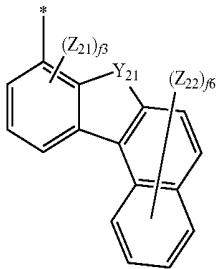
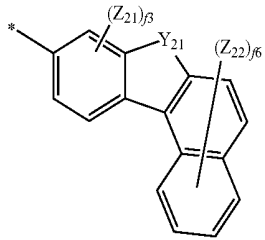
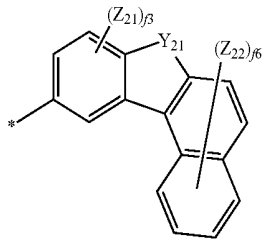
Formula 7-13

Formula 7-14

Formula 7-15

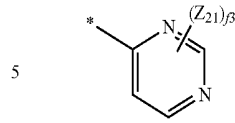
Formula 7-16

31
-continued

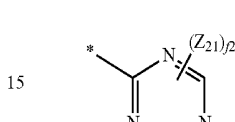


32
-continued

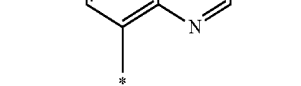
Formula 7-17



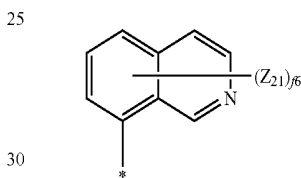
Formula 7-18



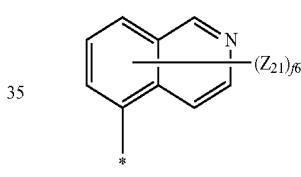
Formula 7-19



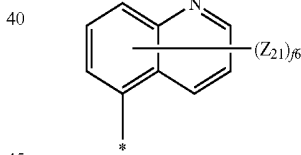
Formula 7-20



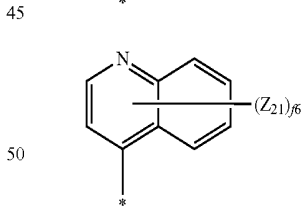
Formula 7-21



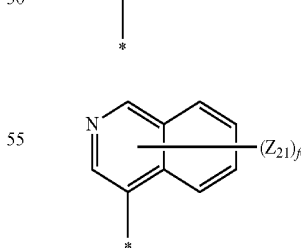
Formula 7-22



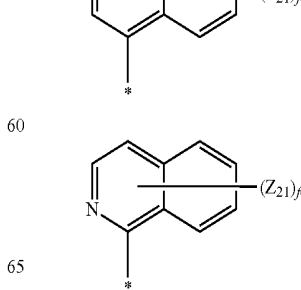
Formula 7-23



Formula 7-24



Formula 7-25



50

Formula 7-26

Formula 7-27

Formula 7-28

Formula 7-29

Formula 7-30

Formula 7-31

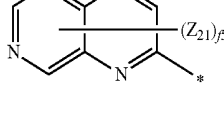
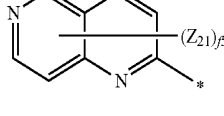
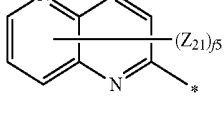
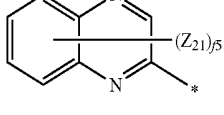
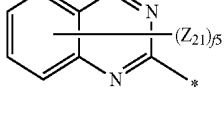
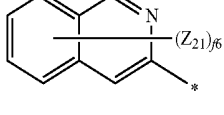
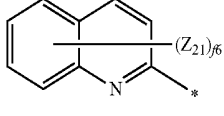
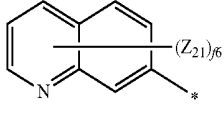
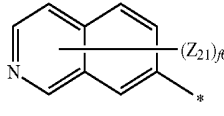
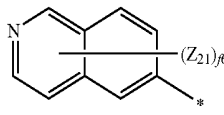
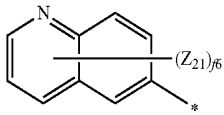
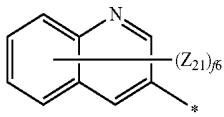
Formula 7-32

Formula 7-33

Formula 7-34

Formula 7-35

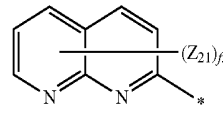
-continued



-continued

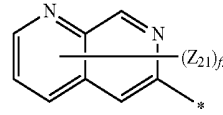
Formula 7-36

5



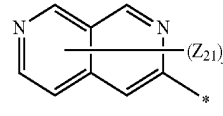
Formula 7-37

10



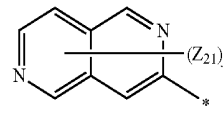
Formula 7-38

15



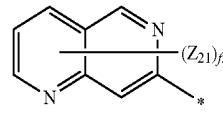
Formula 7-39

20



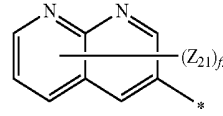
Formula 7-40

25



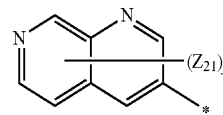
Formula 7-41

30



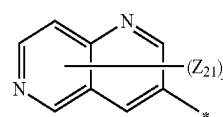
Formula 7-42

35



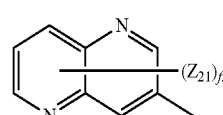
Formula 7-43

40



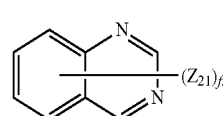
Formula 7-44

45



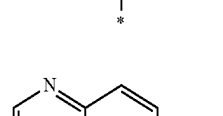
Formula 7-45

50



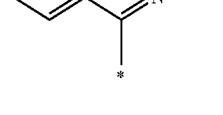
Formula 7-47

55



Formula 7-48

60



65

Formula 7-49

Formula 7-50

Formula 7-51

Formula 7-52

Formula 7-53

Formula 7-54

Formula 7-55

Formula 7-56

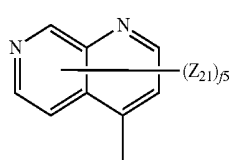
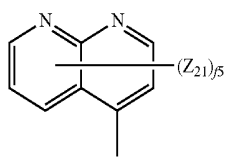
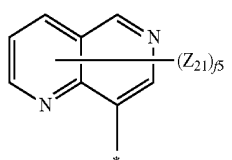
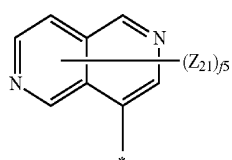
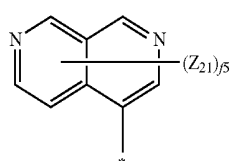
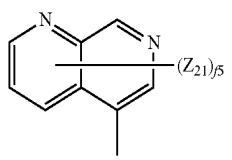
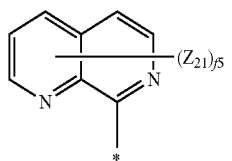
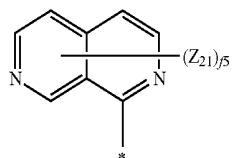
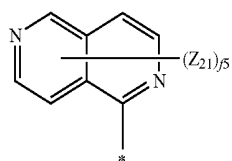
Formula 7-57

Formula 7-58

Formula 7-59

35

-continued

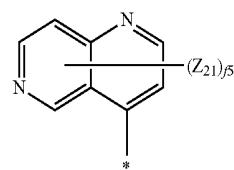


36

-continued

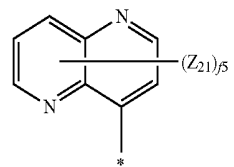
Formula 7-60

5



Formula 7-61

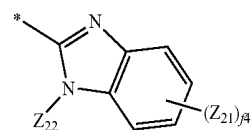
10



15

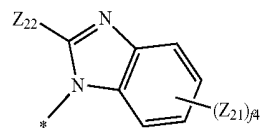
Formula 7-62

20



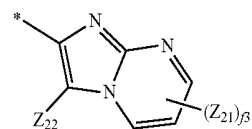
Formula 7-62

25



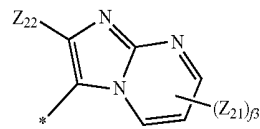
Formula 7-63

30



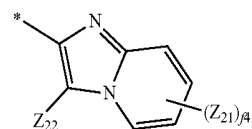
Formula 7-64

35



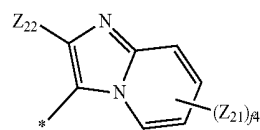
Formula 7-65

40



Formula 7-66

45



50

Formula 7-67

55

wherein, in Formulae 7-1 to 7-75,

Y_{21} may be O, S, C(Z_{23})(Z_{24}), N(Z_{25}), or Si(Z_{26})(Z_{27}),

Z_{21} to Z_{27} 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 phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a triazinyl group, and —Si(Q_{31})(Q_{32})(Q_{33}),

Formula 7-68

Formula 7-69

Formula 7-70

Formula 7-71

Formula 7-72

Formula 7-73

Formula 7-74

Formula 7-75

37

wherein 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,

f2 may be an integer selected from 1 and 2,

f3 may be an integer from 1 to 3,

f4 may be an integer from 1 to 4,

f5 may be an integer from 1 to 5,

f6 may be an integer from 1 to 6,

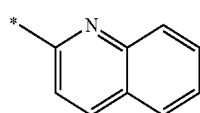
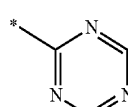
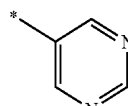
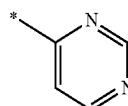
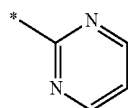
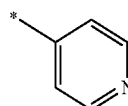
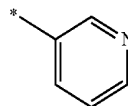
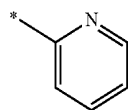
f7 may be an integer from 1 to 7,

f8 may be an integer from 1 to 8,

f9 may be an integer from 1 to 9, and

* indicates a binding site to an adjacent atom.

In one embodiment, in the above formulae, R_1 to R_6 , R_{12} , and R_{13} 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_{10} alkyl group, a C_1 - C_{10} alkoxy group, groups represented by Formulae 8-1 to 8-159, and $-Si(Q_1)(Q_2)(Q_3)$, wherein Q_1 to Q_3 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 are not limited thereto:



Formula 8-1

Formula 8-2

Formula 8-3

Formula 8-4

Formula 8-5

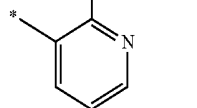
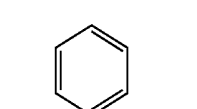
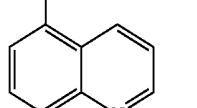
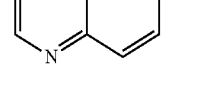
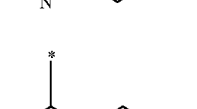
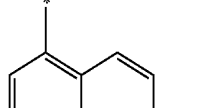
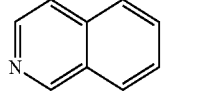
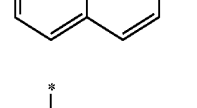
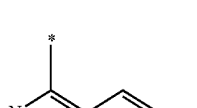
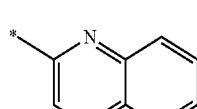
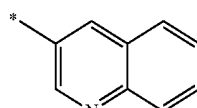
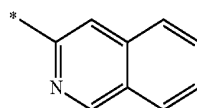
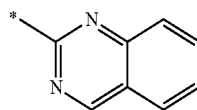
Formula 8-6

Formula 8-7

Formula 8-8

38

-continued



Formula 8-9

Formula 8-10

Formula 8-11

Formula 8-12

Formula 8-13

Formula 8-14

Formula 8-15

Formula 8-16

Formula 8-17

Formula 8-18

5

10

15

20

25

30

35

40

45

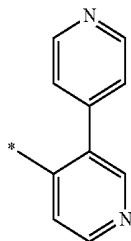
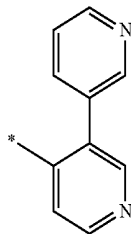
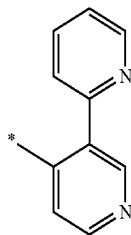
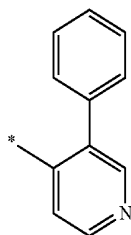
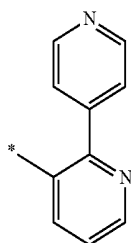
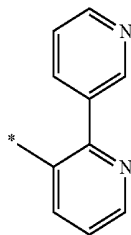
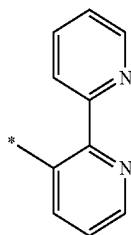
50

55

60

65

39
-continued



Formula 8-19

5

10

Formula 8-20

15

Formula 8-21

20

25

Formula 8-22

30

35

Formula 8-23

40

45

Formula 8-24

50

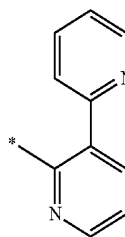
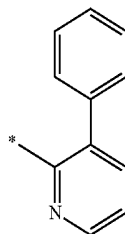
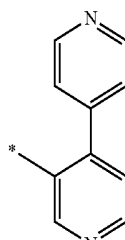
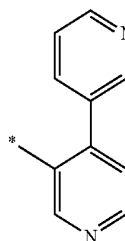
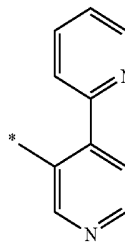
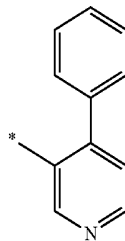
55

Formula 8-25

60

65

40
-continued



Formula 8-26

Formula 8-27

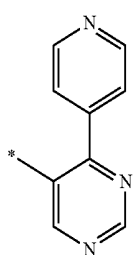
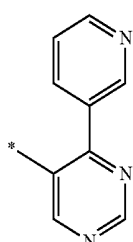
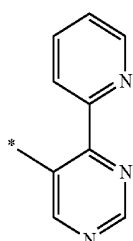
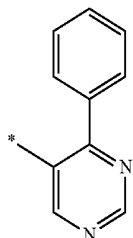
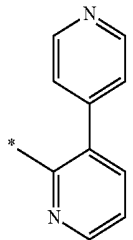
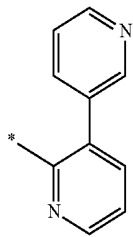
Formula 8-28

Formula 8-29

Formula 8-30

Formula 8-31

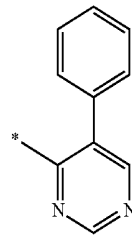
41
-continued



42
-continued

Formula 8-32

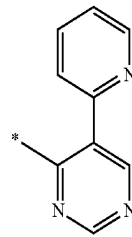
5



10

Formula 8-33

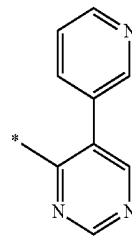
15



20

Formula 8-34

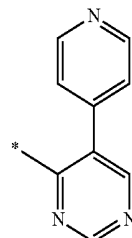
25



30

Formula 8-35

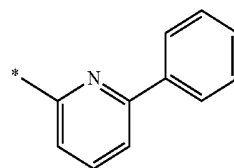
35



40

Formula 8-36

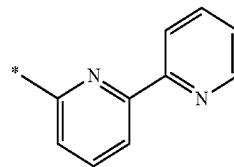
45



50

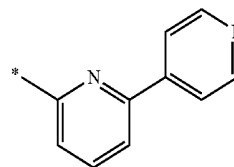
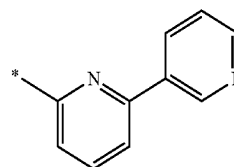
Formula 8-37

60



55

65



Formula 8-38

Formula 8-39

Formula 8-40

Formula 8-41

Formula 8-42

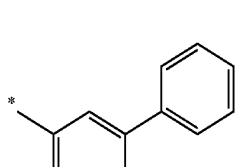
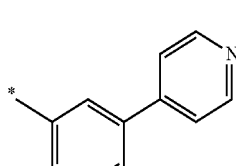
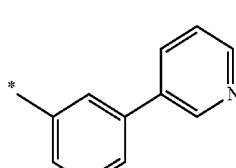
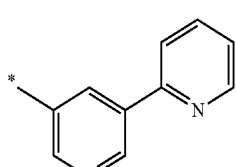
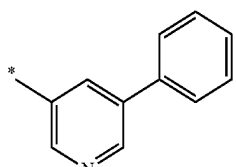
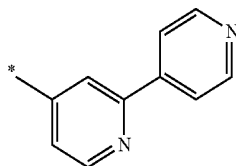
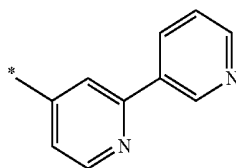
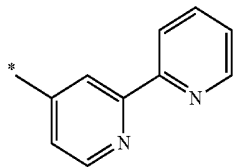
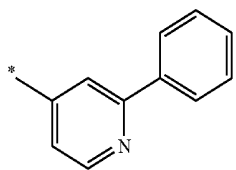
Formula 8-43

Formula 8-44

Formula 8-45

43

-continued

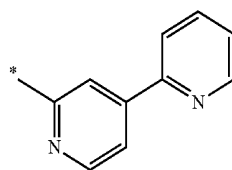


44

-continued

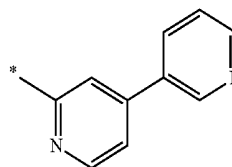
Formula 8-46

5



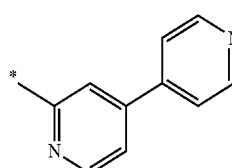
Formula 8-47

10



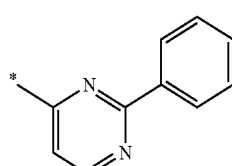
Formula 8-48

15



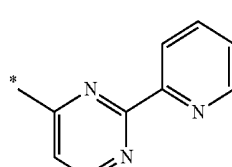
Formula 8-49

25



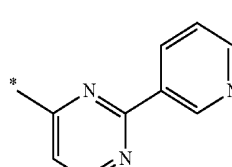
Formula 8-50

30



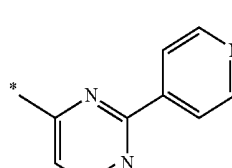
Formula 8-51

35



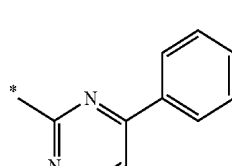
Formula 8-52

45



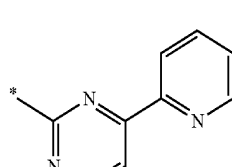
Formula 8-53

55



Formula 8-54

60



Formula 8-55

Formula 8-56

Formula 8-57

Formula 8-58

Formula 8-59

Formula 8-60

Formula 8-61

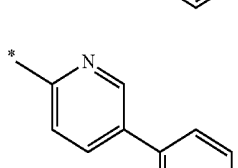
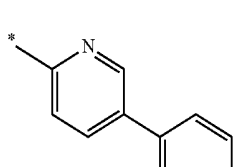
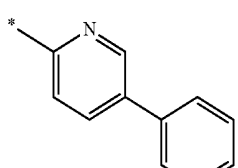
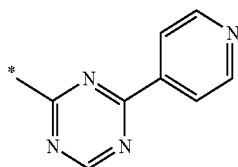
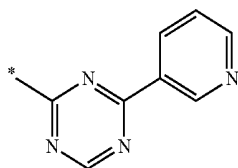
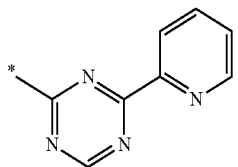
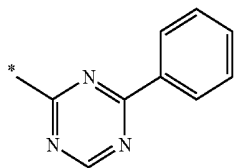
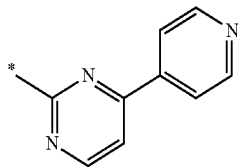
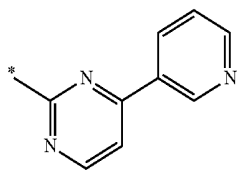
Formula 8-62

Formula 8-63

65

45

-continued

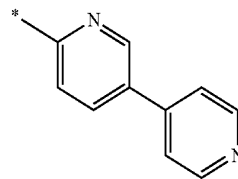


46

-continued

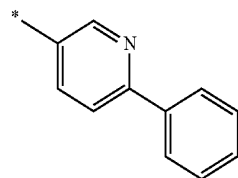
Formula 8-64

5



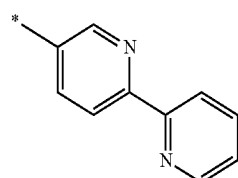
Formula 8-65

10



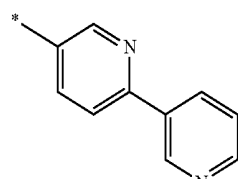
Formula 8-66

15



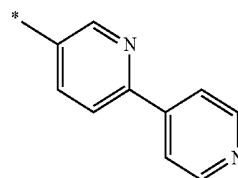
Formula 8-67

25



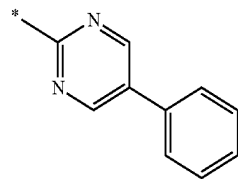
Formula 8-68

30



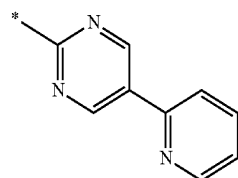
Formula 8-69

40



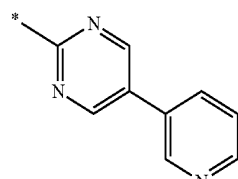
Formula 8-70

45



Formula 8-71

55



Formula 8-72

60

65

Formula 8-73

Formula 8-74

Formula 8-75

Formula 8-76

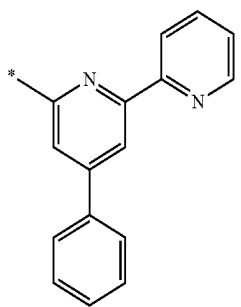
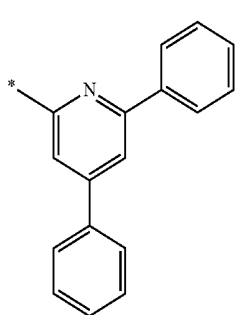
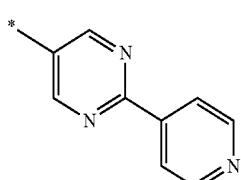
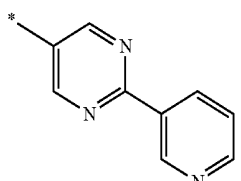
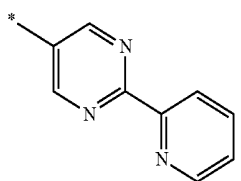
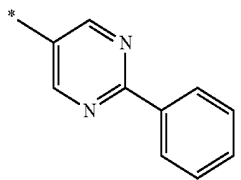
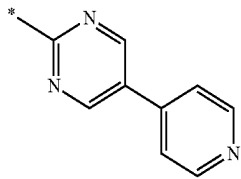
Formula 8-77

Formula 8-78

Formula 8-79

Formula 8-80

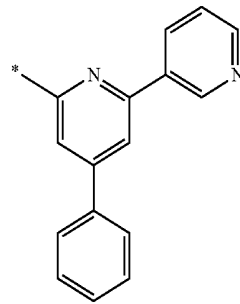
47
-continued



48
-continued

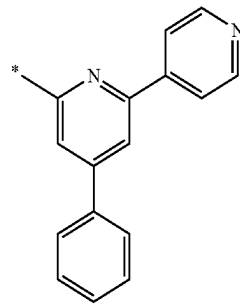
Formula 8-81

5



Formula 8-82 10

15

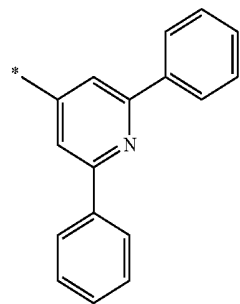


Formula 8-83

20

Formula 8-84 25

30

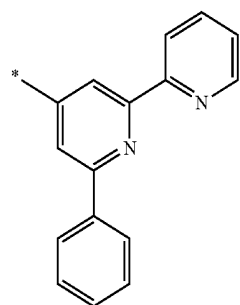


Formula 8-85

35

Formula 8-86 40

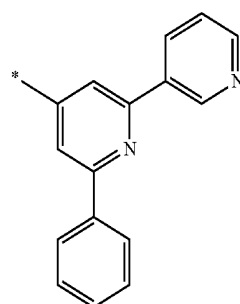
45



50

Formula 8-87 55

60



65

Formula 8-88

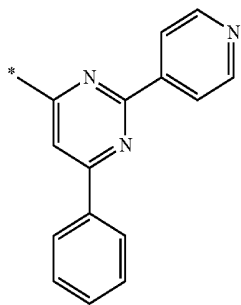
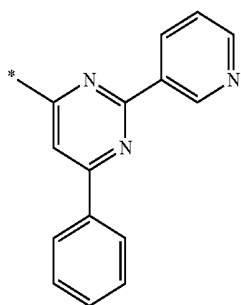
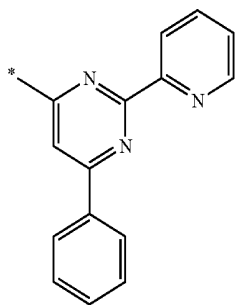
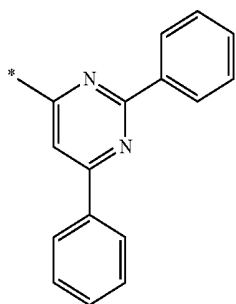
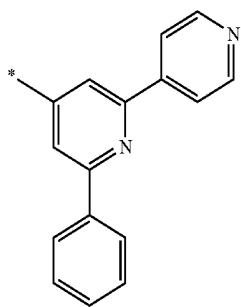
Formula 8-89

Formula 8-90

Formula 8-91

Formula 8-92

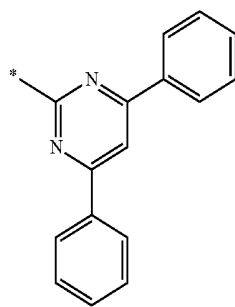
49
-continued



50
-continued

Formula 8-93

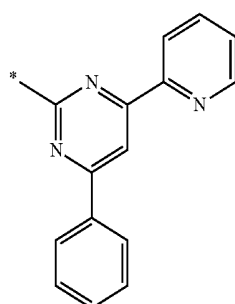
5



10

Formula 8-94

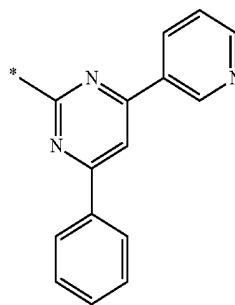
15



20

Formula 8-95

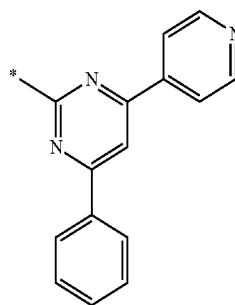
25



35

Formula 8-96

40

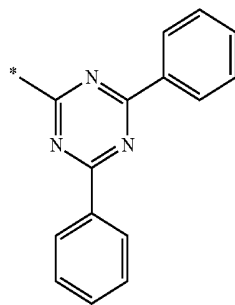


45

50

Formula 8-97

55



60

65

Formula 8-98

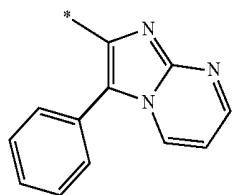
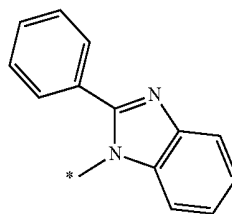
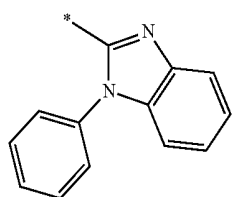
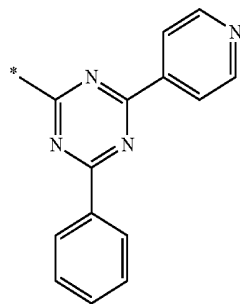
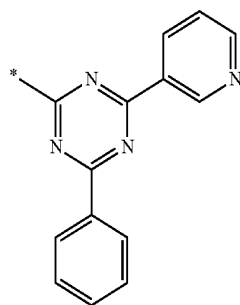
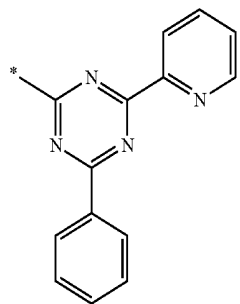
Formula 8-99

Formula 8-100

Formula 8-101

Formula 8-102

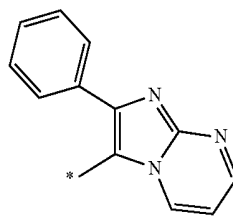
51
-continued



52
-continued

Formula 8-103

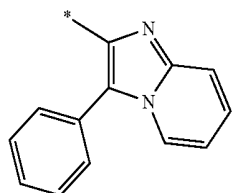
5



10

Formula 8-104

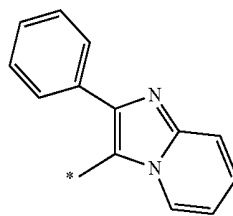
15



20

Formula 8-105

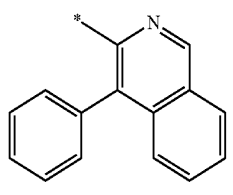
25



30

Formula 8-106

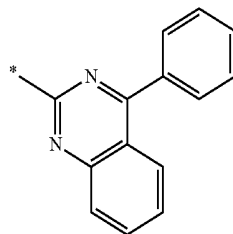
40



45

Formula 8-107

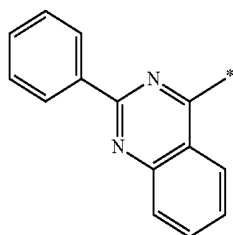
50



55

Formula 8-108

60



65

Formula 8-109

Formula 8-110

Formula 8-111

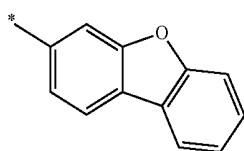
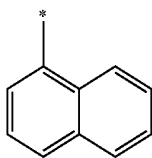
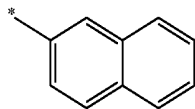
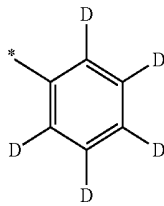
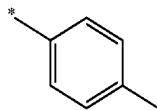
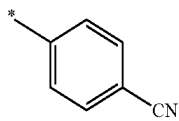
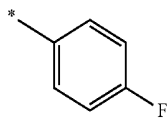
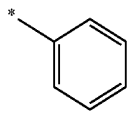
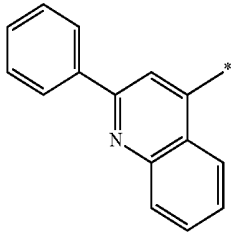
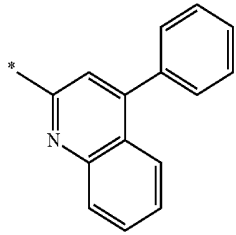
Formula 8-112

Formula 8-113

Formula 8-114

Formula 8-115

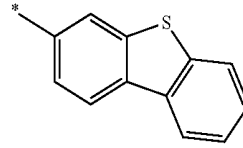
53
-continued



54
-continued

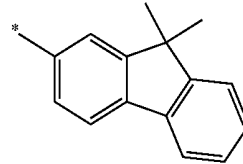
Formula 8-116

5



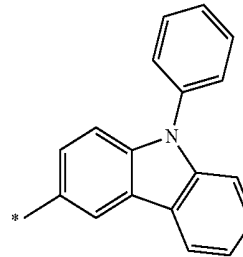
Formula 8-117

10



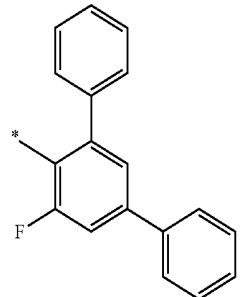
Formula 8-118

20



Formula 8-119

25



Formula 8-120

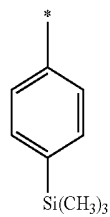
30

Formula 8-121

35

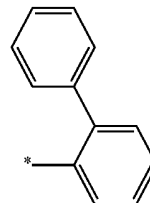
Formula 8-122

40



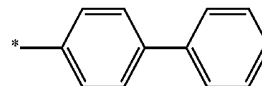
Formula 8-123

50



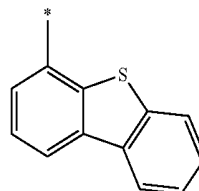
Formula 8-124

55



Formula 8-125

60



65

Formula 8-126

Formula 8-127

Formula 8-128

Formula 8-129

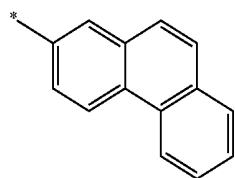
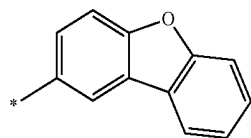
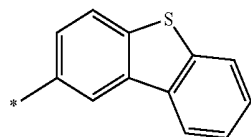
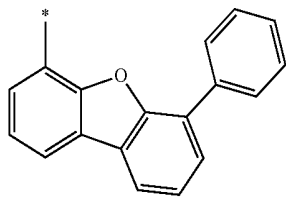
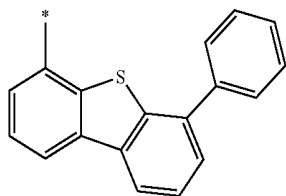
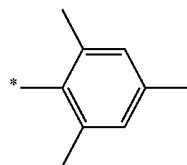
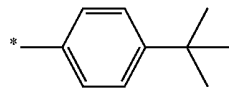
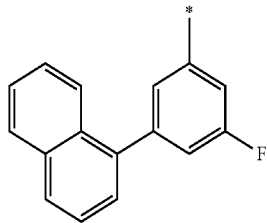
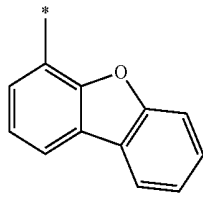
Formula 8-130

Formula 8-131

Formula 8-132

Formula 8-133

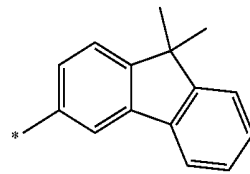
55
-continued



56
-continued

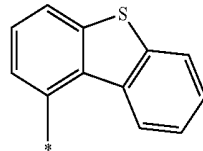
Formula 8-134

5



Formula 8-135

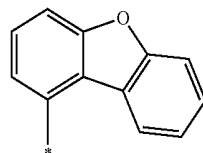
10



15

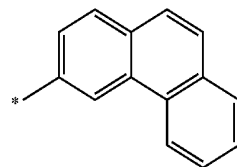
Formula 8-136

20



Formula 8-137

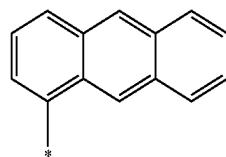
25



30

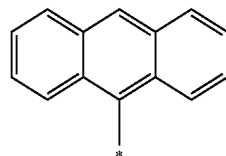
Formula 8-138

35



Formula 8-139

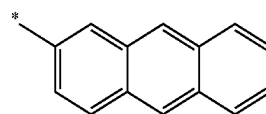
40



45

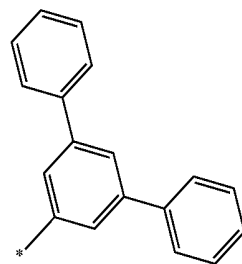
Formula 8-140

50



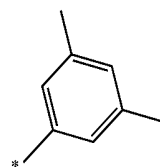
Formula 8-141

55



Formula 8-142

60



65

Formula 8-143

Formula 8-144

Formula 8-145

Formula 8-146

Formula 8-147

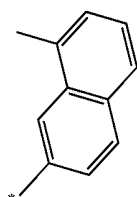
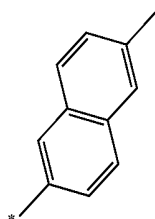
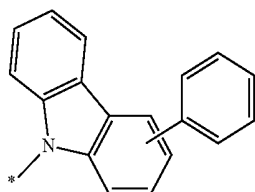
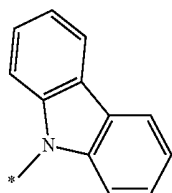
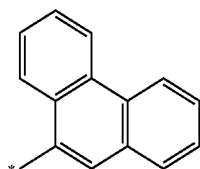
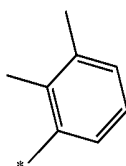
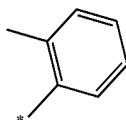
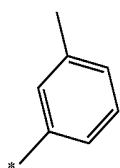
Formula 8-148

Formula 8-149

Formula 8-150

Formula 8-151

-continued



wherein, in Formulae 8-1 to 8-159, * indicates a binding site to an adjacent atom.

In one or more embodiments, in the above formulae, R_3 to R_6 and R_{13} may each independently be selected from hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl

Formula 8-152

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 naphthyl group, and —Si(Q_1)(Q_2)(Q_3),

5 R_{11} , R_2 , and R_{12} 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

Formula 8-153

10 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

Formula 8-154

15 non-aromatic condensed heteropolycyclic group (e.g., groups represented by Formulae 7-1 to 7-75 or groups represented by Formulae 8-1 to 8-159).

In one or more embodiments, in the above formulae,

R_3 to R_6 and R_{13} may be hydrogen,

Formula 8-155

20 R_1 , R_2 , and R_{12} may each independently be selected from groups represented by Formulae 7-1 to 7-75 (e.g., groups represented by Formulae 8-1 to 8-159), but embodiments are not limited thereto.

25 In the above formulae, b1, b2, b5, b6, and b12 may each independently be an integer from 0 to 4; b3 and b4 may each independently be an integer from 0 to 6; and b13 may be 0, 1, or 2.

Formula 8-156

30 In some embodiments, in the above formulae, b1, b2, and b12 may each independently be 0, 1, or 2. In some embodiments, in the above formulae, b1, b2, and b12 may each independently be 1 or 2.

In some embodiments, in the above formulae, b1, b2, and b12 may be 1, but embodiments are not limited thereto.

Formula 8-157

35 In some embodiments, in the above formulae, b3 to b6 and b13 may each independently be 0, 1, or 2. In some embodiments, in the above formulae, b3 to b6 and b13 may each independently be 0 or 1, but embodiments are not limited thereto.

40 In Formula 1, c1 and c2 may each independently be an integer from 0 to 4, and a sum of c1 and c2 may be 1 or greater. That is, at least one selected from a group represented by $^*-(L_1)_{a1}-(R_1)_{b1}$ and a group represented by $^*-(L_2)_{a2}-(R_2)_{b2}$ may be essentially present in Formula 1.

Formula 8-158

In one embodiment, in Formula 1, a sum of c1 and c2 may be 1 or 2.

45 In one or more embodiments, in Formula 1,

c1 may be 1, and c2 may be 0;

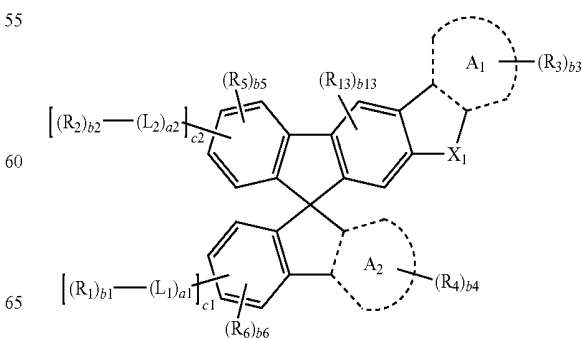
c1 may be 1, and c2 may be 1; or

c1 may be 0, and c2 may be 1, but embodiments are not limited thereto.

50 In one embodiment, the condensed-cyclic compound may be represented by one of Formulae 1A to 1E:

Formula 8-159

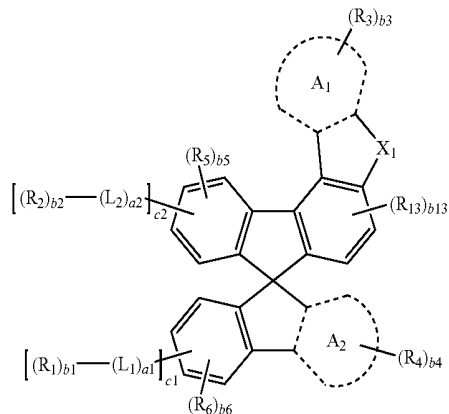
Formula 1A



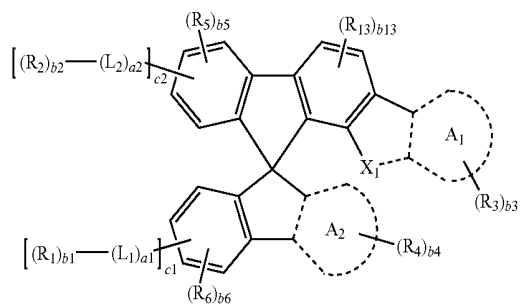
59

-continued

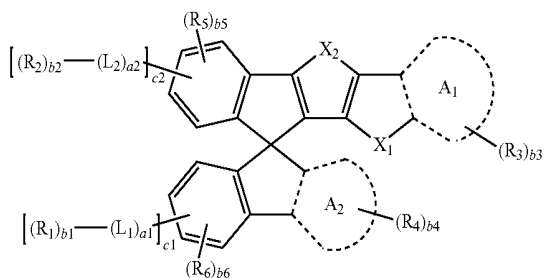
Formula 1B



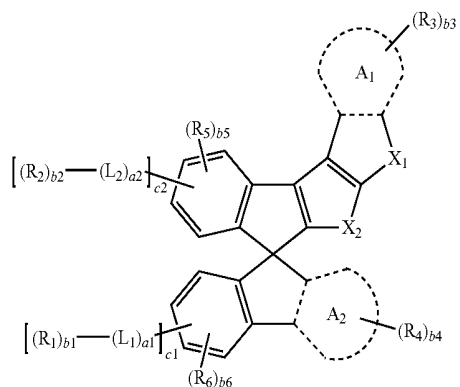
Formula 1C



Formula 1D



Formula 1E



wherein, in Formulae 1A to 1E, ring A₁, ring A₂, X₁, X₂, L₁, L₂, a₁, a₂, L₁₁, L₁₂, a₁₁, a₁₂, R₁ to R₆, R₁₁ to R₁₃, b₁ to b₆, b₁₁ to b₁₃, c₁, and c₂ may be the same as those described herein.

In some embodiments, in Formulae 1A and 1E,

ring A₁ may be a benzene group or a naphthalene group, ring A₂ may be selected from a pyridine group, a quinoline group, and an isoquinoline group,

X₁ may be N-[(L₁₁)_{a11}-(R₁₁)_{b11}],

X₂ may be O or S,

60

L₁ and L₂ may each independently be selected from groups represented by Formulae 3-8, 3-9, 3-25, and 3-35 to 3-41 (e.g., groups represented by Formulae 4-11, 4-13, 4-27 and 4-29 to 4-35),

a₁ and a₂ may each independently be 0, 1, or 2,

L₁₁ may be selected from groups represented by Formulae 3-1 to 3-9 (provided that Y₁ in Formulae 3-3 and 3-4 may be C(Z₃)(Z₄)), 3-25, and 3-33 to 3-41 (e.g., groups represented by Formulae 4-1, 4-3, 4-5, 4-7 to 4-13, 4-17, and 4-24 to 4-35),

a₁₁ may be 0, 1, or 2,

R₁₁ may be selected from groups represented by Formulae 5-1 to 5-19 (e.g., groups represented by Formulae 6-1 to 6-27),

b₁₁ may be 1 or 2,

R₁ and R₂ may each independently be selected from groups represented by Formulae 7-1 to 7-75 (e.g., groups represented by Formulae 8-1 to 8-159),

b₁ and b₂ may each independently be 1 or 2,

R₃ to R₆ and R₁₃ 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, and —Si(Q₁)(Q₂)(Q₃),

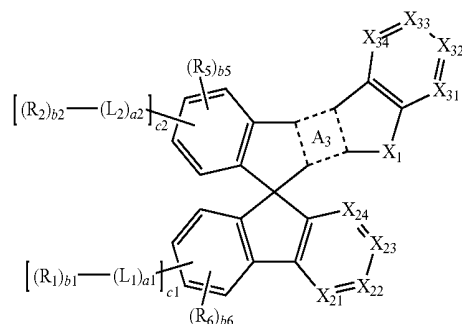
wherein Q₁ to Q₃ may each independently be selected from a C₁-C₁₀ alkyl group, a C₁-C₁₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group,

b₃ to b₆ and b₁₃ may each independently be 0, 1, or 2, and

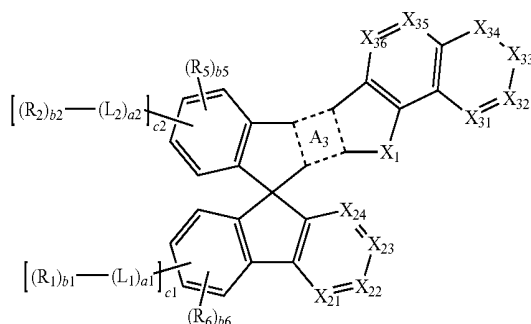
c₁ may be 1, and c₂ may be 0; c₁ may be 1, and c₂ may be 1; or c₁ may be 0, and c₂ may be 1, but embodiments are not limited thereto.

In one or more embodiments, the condensed-cyclic compound may be represented by one of Formulae 1-1 to 1-7:

Formula 1-1



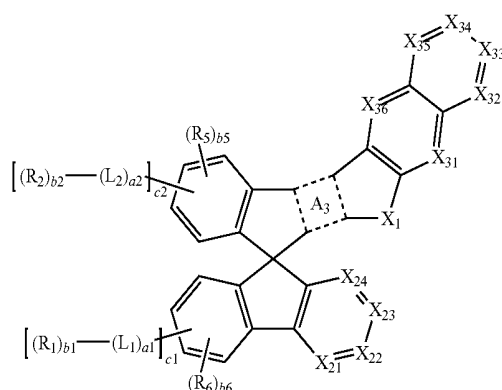
Formula 1-2



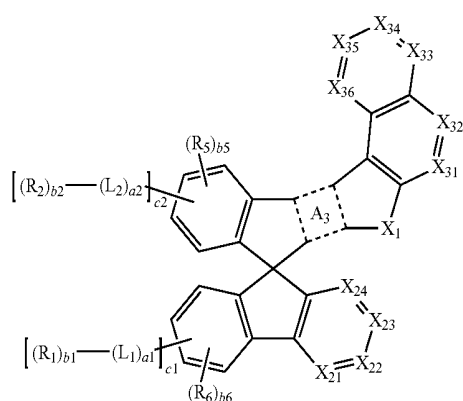
61

-continued

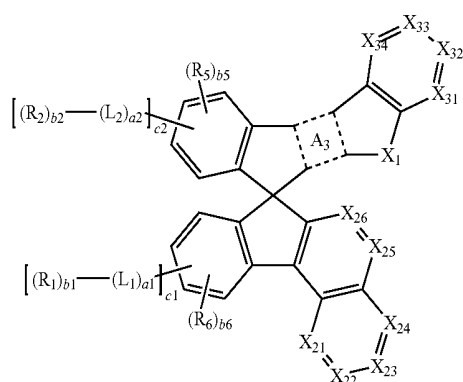
Formula 1-3



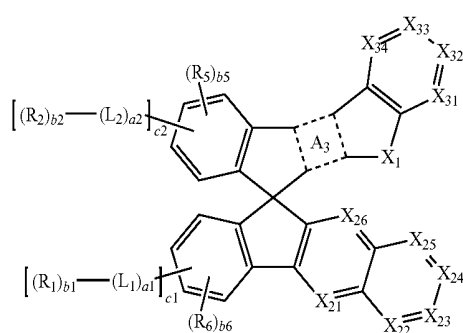
Formula 1-4



Formula 1-5



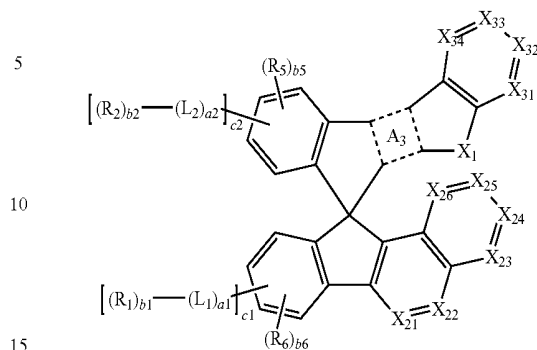
Formula 1-6



62

-continued

Formula 1-7



wherein, in Formulae 1-1 to 1-7,

ring A₃, X₁, X₂, L₁, L₂, a₁, a₂, L₁₁, L₁₂, a₁₁, a₁₂, R₁, R₂, R₅, R₆, R₁₁ to R₁₃, b₁, b₂, b₅, b₆, b₁₁ to b₁₃, c₁, and c₂ may be the same as those described herein,

X₂₁ may be N or G(R₂₁), X₂₂ may be N or G(R₂₂), X₂₃ may be N or G(R₂₃), X₂₄ may be N or G(R₂₄), X₂₅ may be N or G(R₂₅), X₂₆ may be N or G(R₂₆),

at least one selected from X₂₁ to X₂₆ may be N,

X₃₁ may be N or C(R₃₁), X₃₂ may be N or C(R₃₂), X₃₃ may be N or C(R₃₃), X₃₄ may be N or C(R₃₄), X₃₅ may be N or C(R₃₅), X₃₆ may be N or C(R₃₆),

R₂₁ and R₂₆ may each independently be the same as describe herein with reference to R₃, and

R₃₁ and R₃₆ may each independently be the same as describe herein with reference to R₄.

In one embodiment,

in Formula 1-1, one or two of X₂₁ to X₂₄ may be N, and none, one, or two of X₃₁ to X₃₄ may be N,

in Formulae 1-2 to 1-4, one or two of X₂₁ to X₂₄ may be N, and none, one, or two of X₃₁ to X₃₆ may be N,

in Formulae 1-5 to 1-7, one or two of X₂₁ to X₂₆ may be N, and none, one, or two of X₃₁ to X₃₆ may be N.

In one or more embodiments,

in Formulae 1-1 to 1-4, one of X₂₁ to X₂₄ may be N,

in Formulae 1-1 and 1-5 to 1-7, none or one of X₃₁ to X₃₄ may be N,

in Formulae 1-5 to 1-7, one of X₂₁ to X₂₆ may be N, and in Formulae 1-2 to 1-4, none or one of X₃₁ to X₃₆ may be N.

In some embodiments, in Formulae 1-1 to 1-7, ring A₃ may be a group represented by Formula 2A.

In one embodiment, in Formulae 1-1 to 1-7,

X₁ may be N-[(L₁₁)_{a11}-(R₁₁)_{b11}],

X₂ may be O or S,

L₁ and L₂ may each independently be selected from groups represented by Formulae 3-8, 3-9, 3-25, and 3-35 to 3-41 (e.g., groups represented by Formulae 4-11, 4-13, 4-27 and 4-29 to 4-35),

a₁ and a₂ may each independently be 0, 1, or 2,

L₁₁ may be selected from groups represented by Formulae 3-1 to 3-9 (provided that Y₁ in Formulae 3-3 and 3-4 may be C(Z₃)(Z₄)), 3-25, and 3-33 to 3-41 (e.g., groups represented by Formulae 4-1, 4-3, 4-5, 4-7 to 4-13, 4-17, and 4-24 to 4-35),

a₁₁ may be 0, 1, or 2,

R₁₁ may be selected from groups represented by Formulae 5-1 to 5-19 (e.g., groups represented by Formulae 6-1 to 6-27),

63

b11 may be 1 or 2,

R₁ and R₂ may each independently be selected from groups represented by Formulae 7-1 to 7-75 (e.g., groups represented by Formulae 8-1 to 8-159),

b1 and b2 may each independently be 1 or 2,

R₃ to R₆ and R₁₃ 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, and —Si(Q₁)(Q₂)(Q₃),

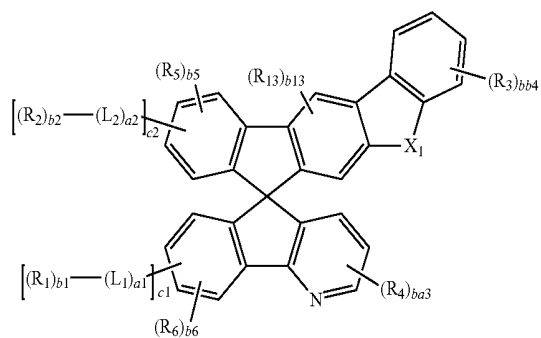
wherein Q₁ to Q₃ may each independently be selected from a C₁-C₁₀ alkyl group, a C₁-C₁₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group,

b3 to b6 and b13 may each independently be 0, 1, or 2, and

c1 may be 1, and c2 may be 0; c1 may be 1, and c2 may be 1; or c1 may be 0, and c2 may be 1, but embodiments are not limited thereto.

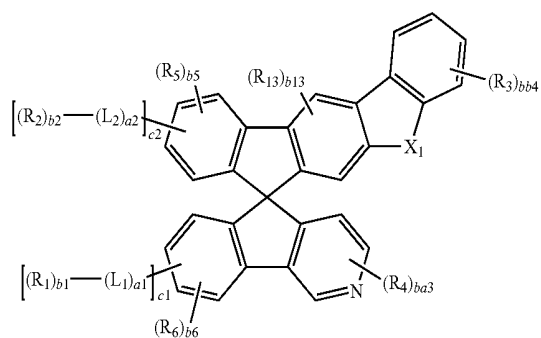
In one or more embodiments, the condensed-cyclic compound may be represented by one of Formulae 1(1) to 1(19), but embodiments are not limited thereto:

Formula 1(1)

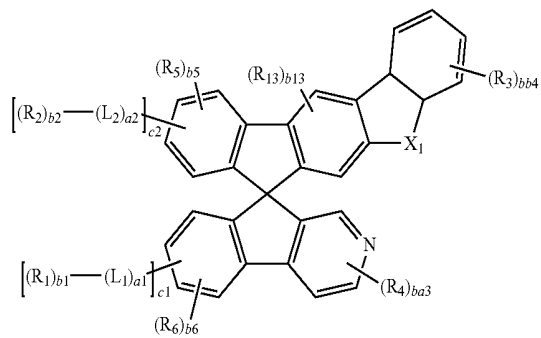


Formula 1(1)

Formula 1(2)



Formula 1(3)

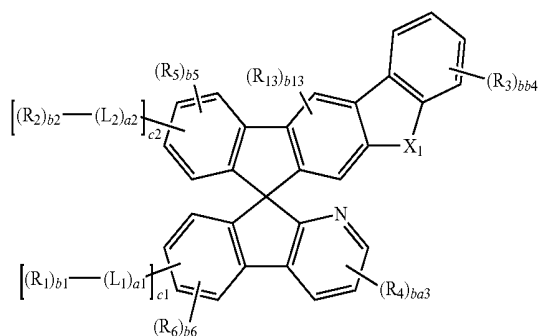


Formula 1(3)

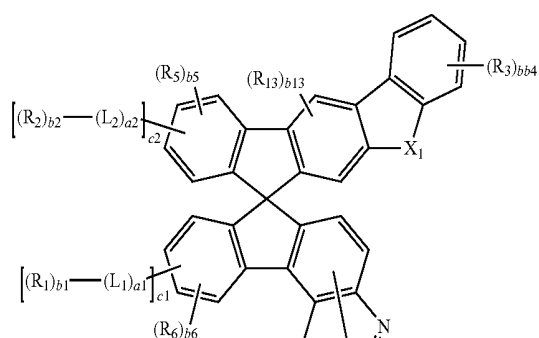
64

-continued

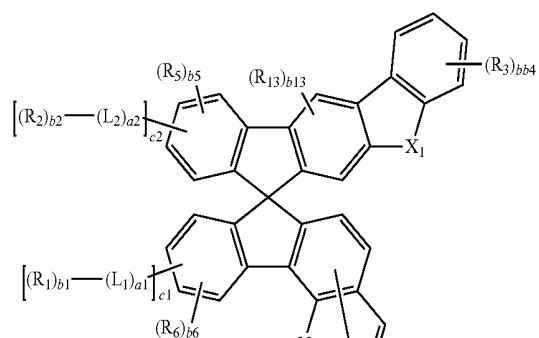
Formula 1(4)



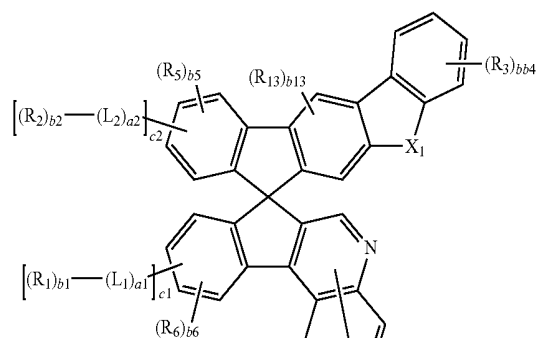
Formula 1(5)



Formula 1(6)



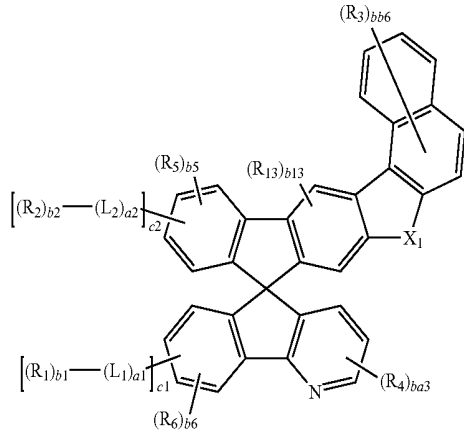
Formula 1(7)



65

65
-continued

Formula 1(8)



5

10

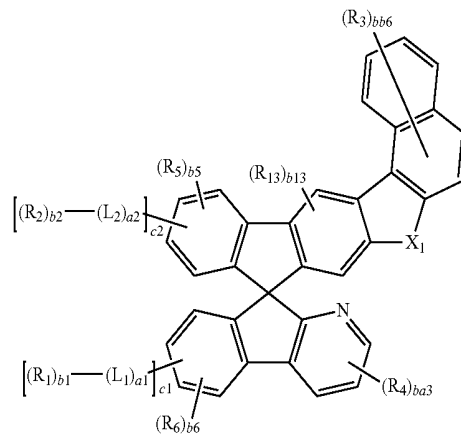
15

20

25

66
-continued

Formula 1(11)



5

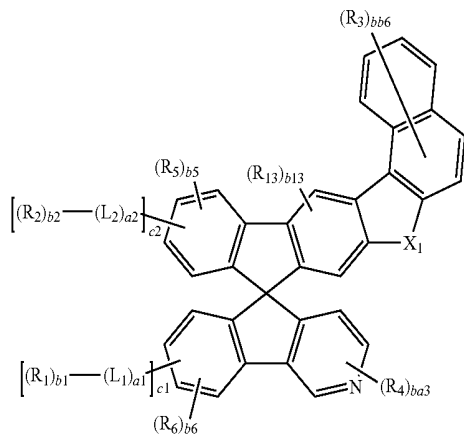
10

15

20

25

Formula 1(9)



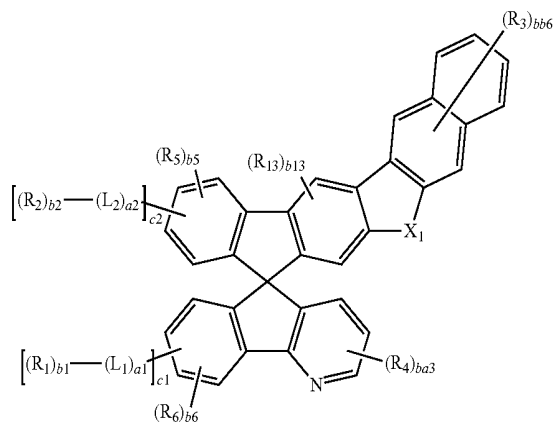
30

35

40

45

Formula 1(12)



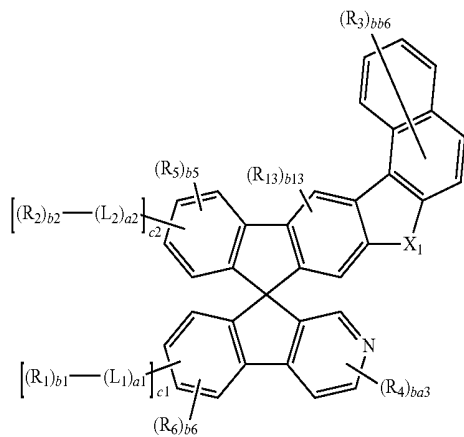
30

35

40

45

Formula 1(10)



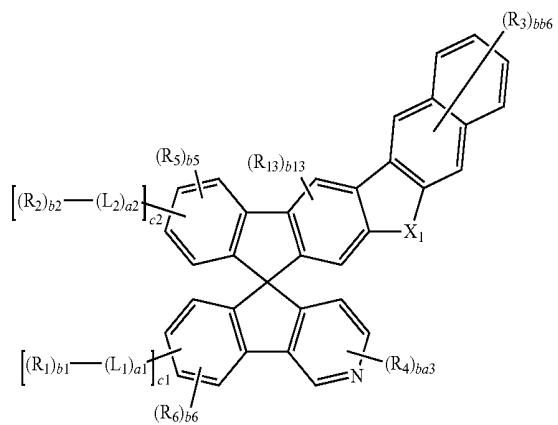
50

55

60

65

Formula 1(13)



55

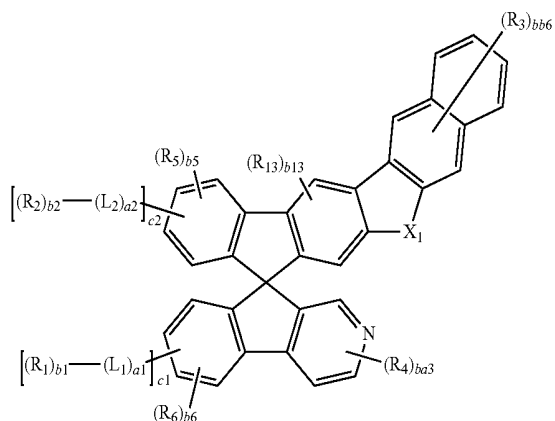
60

65

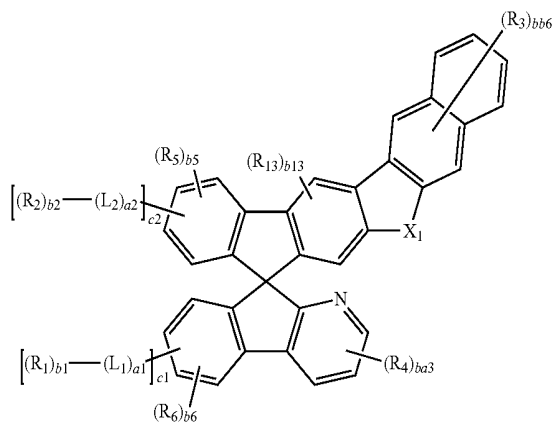
67

-continued

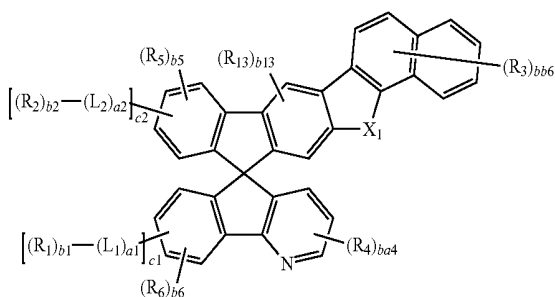
Formula 1(14)



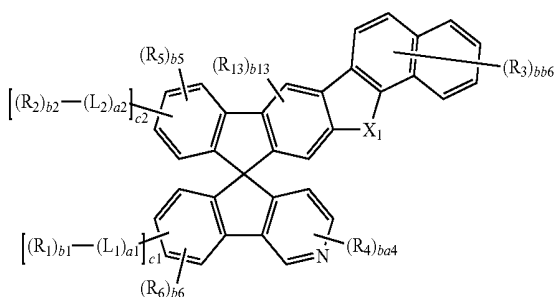
Formula 1(15)



Formula 1(16)



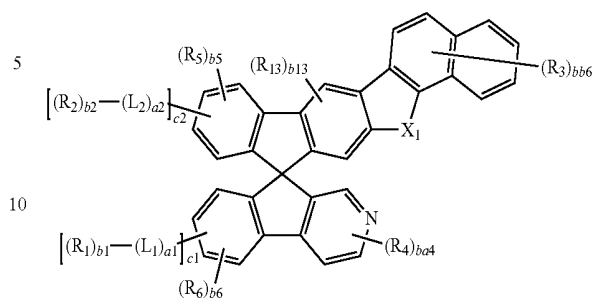
Formula 1(17)



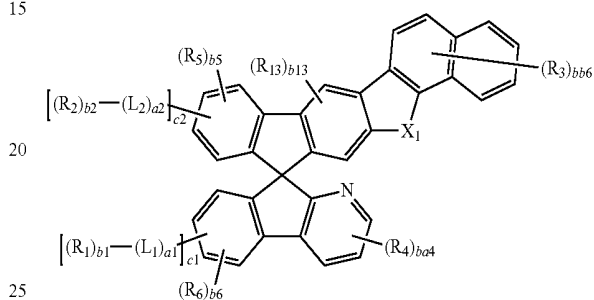
68

-continued

Formula 1(18)



Formula 1(19)



wherein, in Formulae 1(1) to 1(19),

X_1 , L_1 , L_2 , a_1 , a_2 , L_{11} , L_{12} , a_{11} , a_2 , R_1 to R_6 , R_{11} to R_{13} , b_1 to b_6 , b_1 to b_3 ,

c_1 , and c_2 may be the same as those described herein,

ba_3 may be an integer from 0 to 3,

bb_4 may be an integer from 0 to 4,

ba_5 may be an integer from 0 to 5, and

bb_6 may be an integer from 0 to 6.

In one embodiment, in Formulae 1(1) to 1(19),

X_1 may be $N-[(L_{11})_{a11}-(R_{11})_{b11}]$,

X_2 may be O or S,

L_1 and L_2 may each independently be selected from groups represented by Formulae 3-8, 3-9, 3-25, and 3-35 to 3-41 (e.g., groups represented by Formulae 4-11, 4-13, 4-27 and 4-29 to 4-35),

a_1 and a_2 may each independently be 0, 1, or 2,

L_{11} may be selected from groups represented by Formulae 3-1 to 3-9 (provided that Y_1 in Formulae 3-3 and 3-4 may be $C(Z_3)(Z_4)$), 3-25, and 3-33 to 3-41 (e.g., groups represented by Formulae 4-1, 4-3, 4-5, 4-7 to 4-13, 4-17, and 4-24 to 4-35),

a_{11} may be 0, 1, or 2,

R_{11} may be selected from groups represented by Formulae 5-1 to 5-19 (e.g., groups represented by Formulae 6-1 to 6-27),

b_{11} may be 1 or 2,

R_1 and R_2 may each independently be selected from groups represented by Formulae 7-1 to 7-75 (e.g., groups represented by Formulae 8-1 to 8-159),

b_1 and b_2 may each independently be 1 or 2,

R_3 to R_6 and R_{13} 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 phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, and $-Si(Q_1)(Q_2)(Q_3)$,

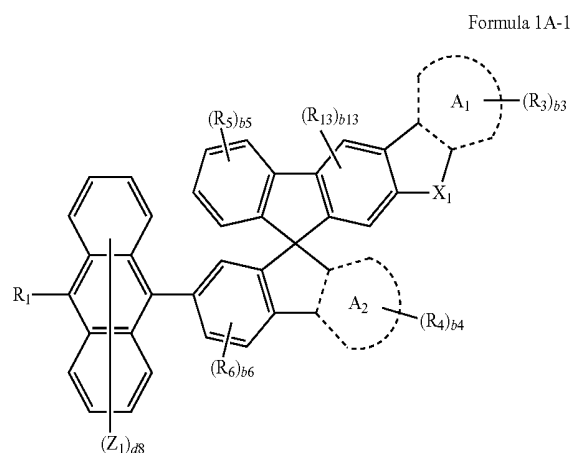
wherein Q_1 to Q_3 may each independently be selected from a C_1 - C_{10} alkyl group, a C_1 - C_{10} alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group,

69

ba3, bb4, ba5, bb6, b5, and b6 may each independently be 0, 1, or 2, and

c1 may be 1, and c2 may be 0; c1 may be 1, and c2 may be 1; or c1 may be 0, and c2 may be 1, but embodiments are not limited thereto.

In one or more embodiments, the condensed-cyclic compound may be represented by one of Formulae 1A-1 to 1A-3, but embodiments are not limited thereto:



wherein, in Formulae 1A-1 to 1A-3, ring A, ring A₂, X₁, L₁₁, L₁₂, a₁₁, a₁₂, R₁, R₃ to R₆, R₁₁ to R₁₃, b₃ to b₆, b₁₁ to b₁₃, c₁, and c₂ may be the same as those described herein.

In some embodiments, in Formulae 1A-1 to 1A-3, ring A₁ may be selected from a benzene group and a naphthalene group, ring A₂ may be selected from a pyridine group, a quinoline group, and an isoquinoline group,

70

X₁ may be N-[(L₁₁)_{a11}-(R₁₁)_{b11}],

L₁₁ may be selected from groups represented by Formulae 3-1 to 3-9 (provided that Y₁ in Formulae 3-3 and 3-4 may be C(Z₃)(Z₄)), 3-25, and 3-33 to 3-41 (e.g., groups represented by Formulae 4-1, 4-3, 4-5, 4-7 to 4-13, 4-17, and 4-24 to 4-35),

a₁₁ may be 0, 1, or 2,

R₁₁ may be selected from groups represented by Formulae 5-1 to 5-19 (e.g., groups represented by Formulae 6-1 to 6-27),

b₁₁ may be 1 or 2,

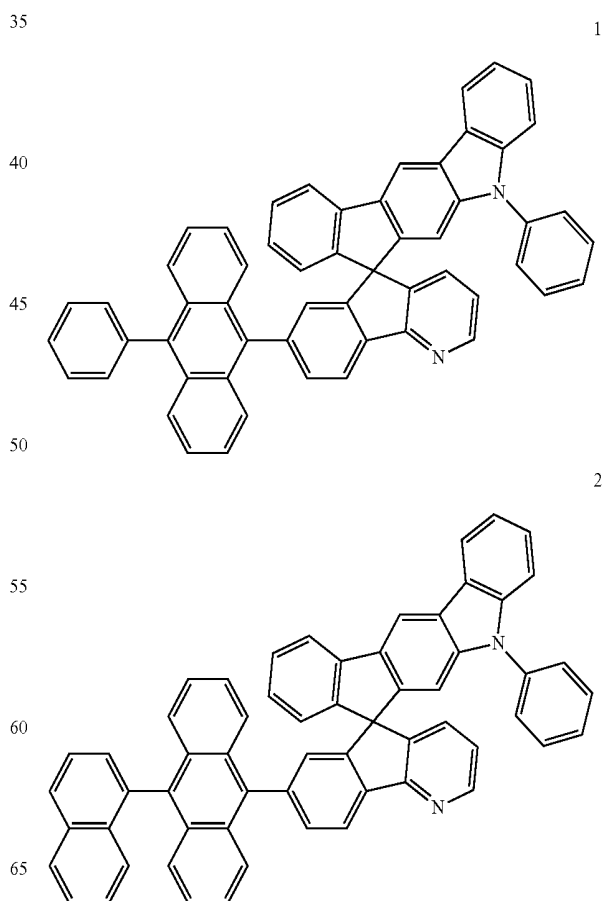
R₁ may be selected from groups represented by Formulae 7-1 to 7-75 (e.g., groups represented by Formulae 8-1 to 8-159),

b₁ may be 1 or 2,

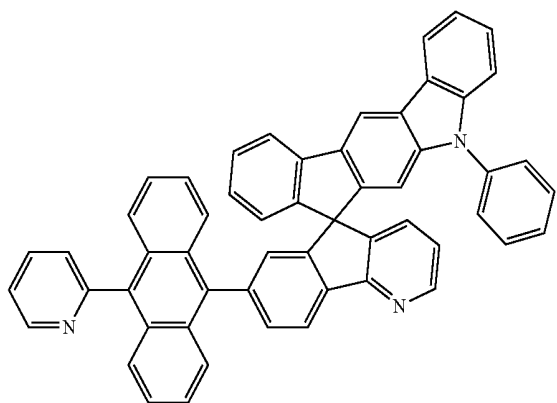
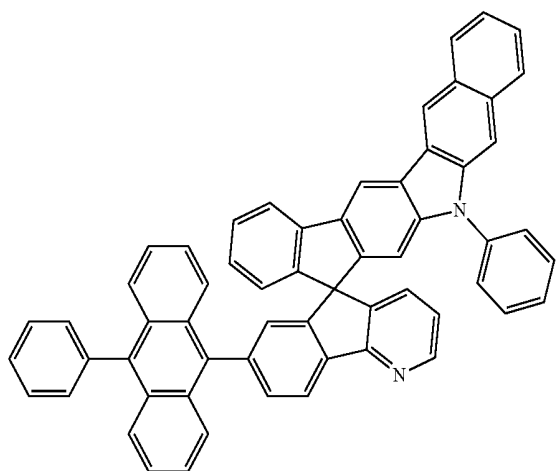
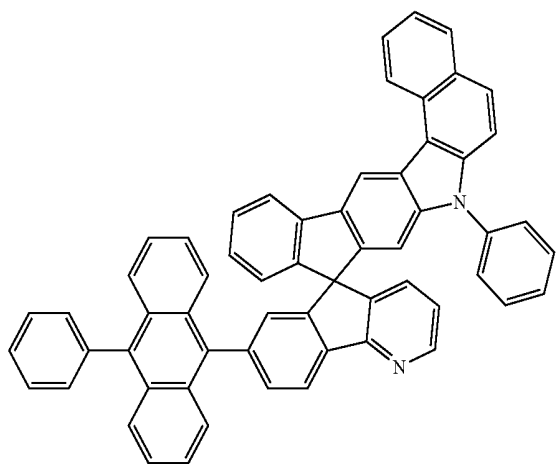
R₃ to R₆, R₁₃, Z₁, and Z₂ may each independently be selected from hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group, and

b₃ to b₆, b₁₃, d₄, d₅, and d₈ may each independently be selected from 0, 1, and 2, but embodiments are not limited thereto:

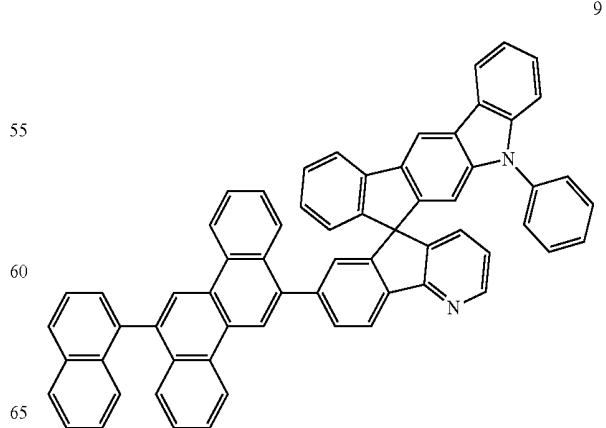
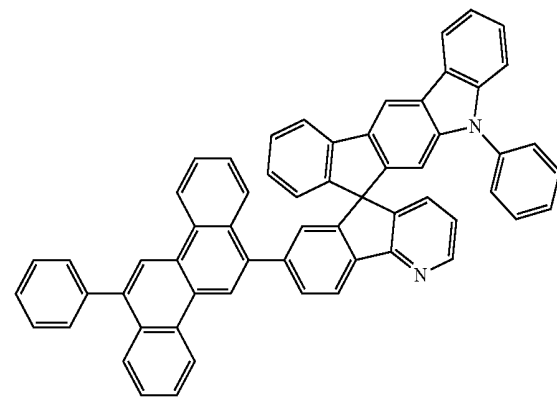
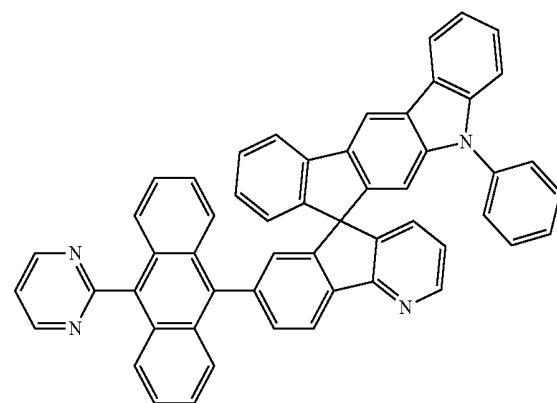
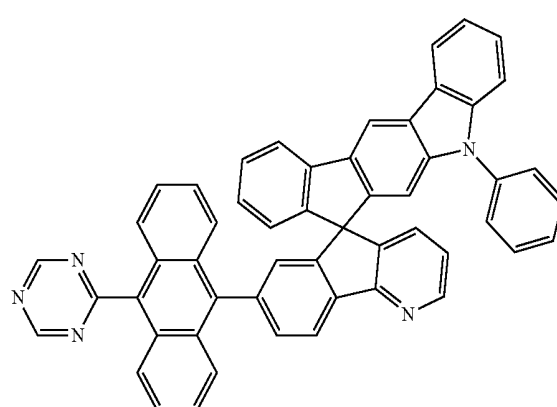
In some embodiments, the condensed-cyclic compound represented by Formula 1 may be represented by one of Compounds 1 to 35, but embodiments are not limited thereto:



71
-continued

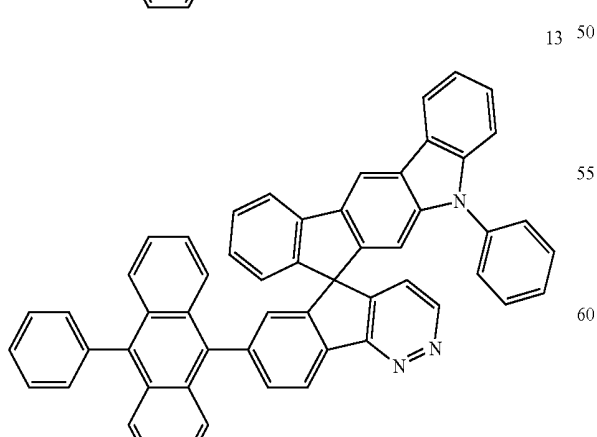
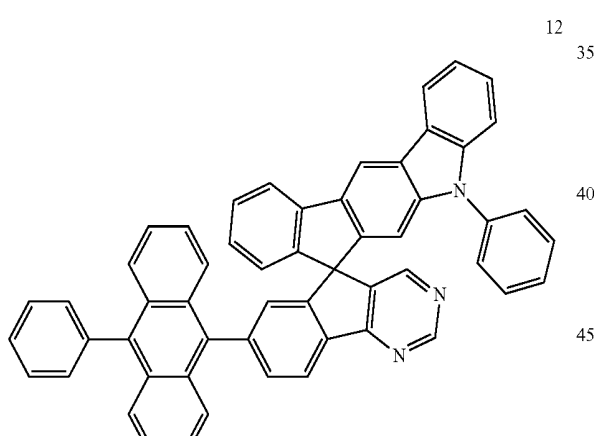
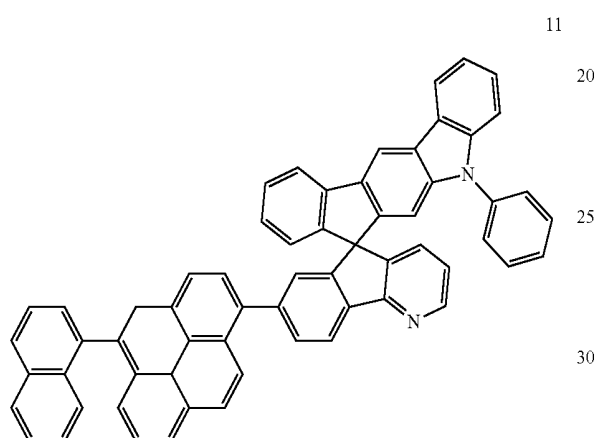
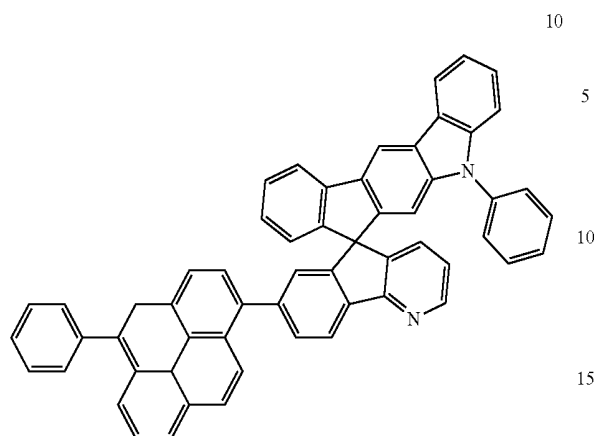


72
-continued



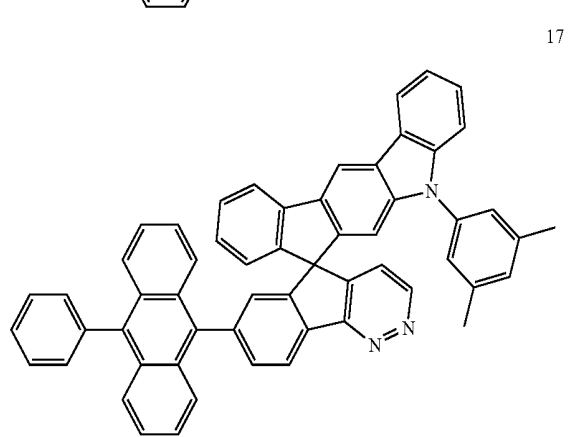
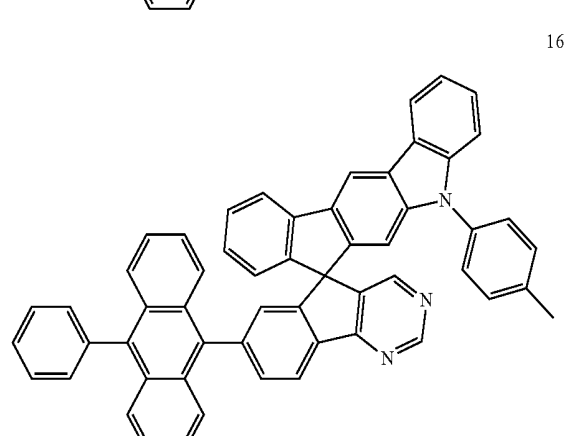
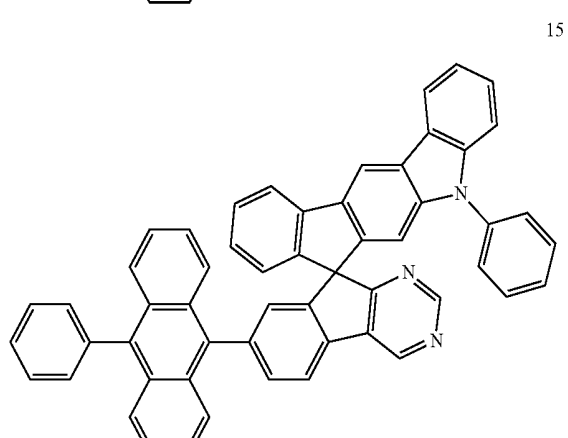
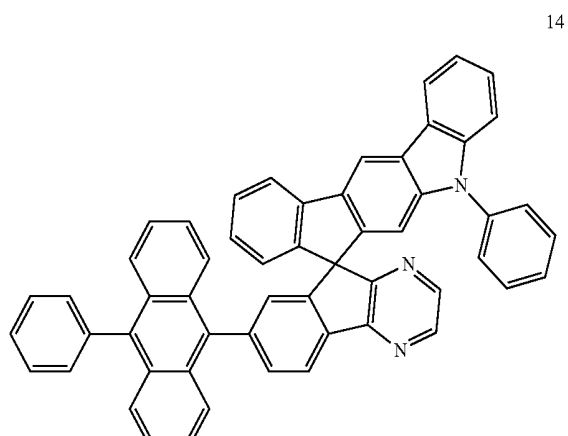
73

-continued



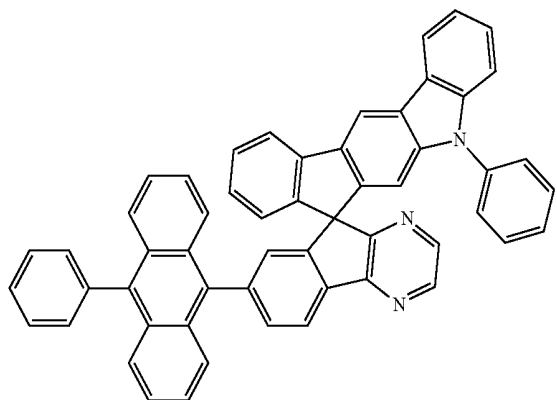
74

-continued



75
-continued

18



5

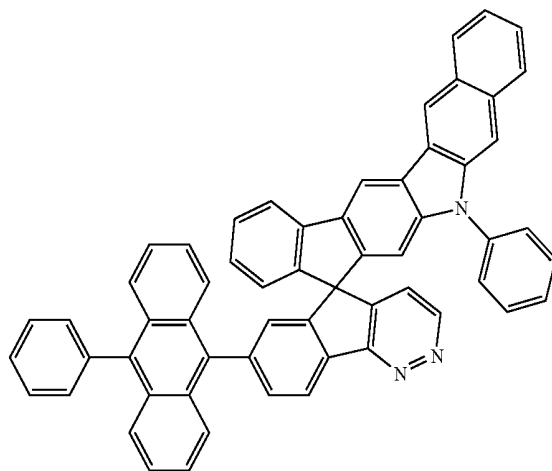
10

15

20

76
-continued

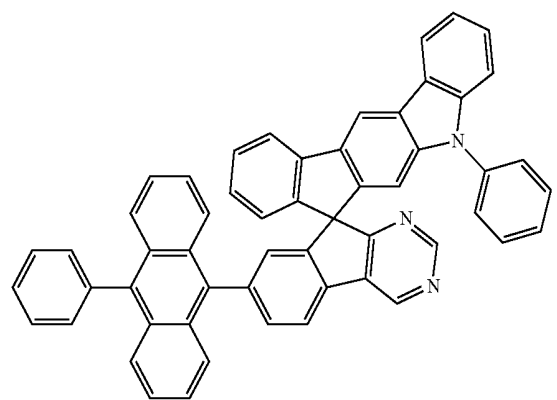
21



19

25

22

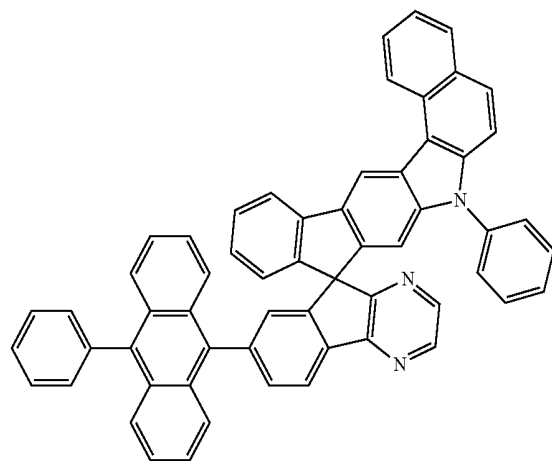


30

35

40

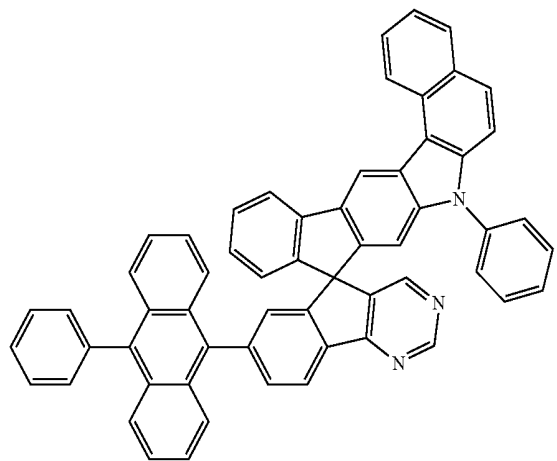
45



20

50

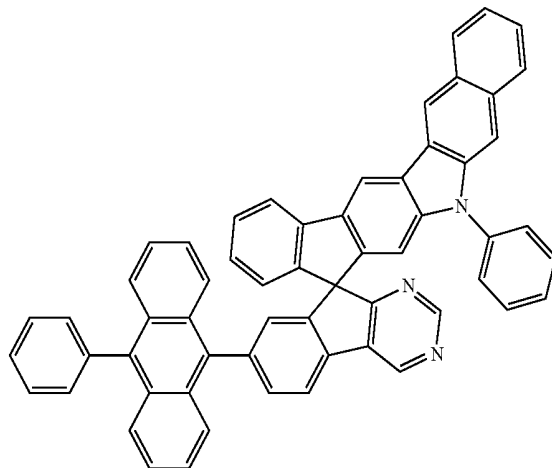
23



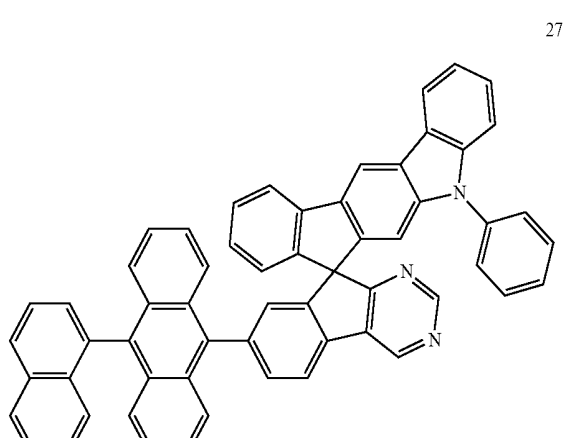
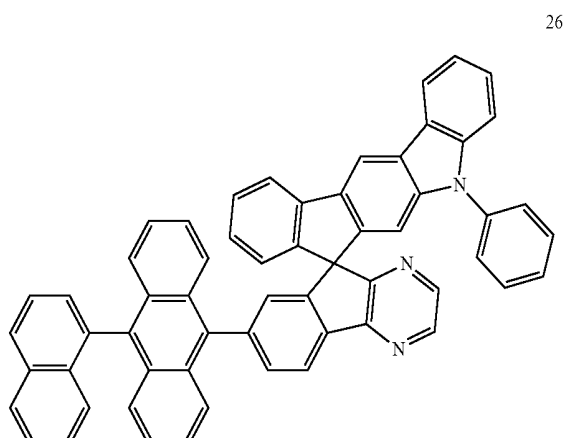
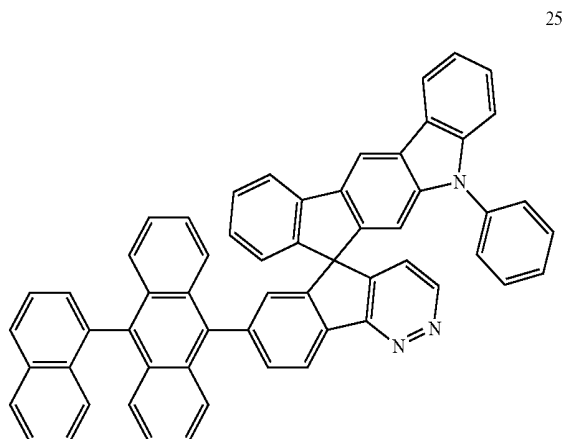
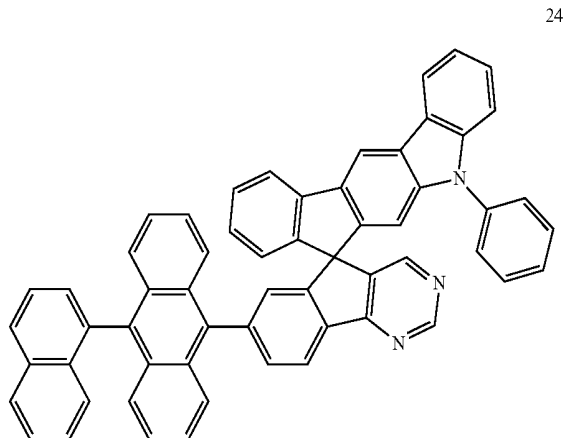
55

60

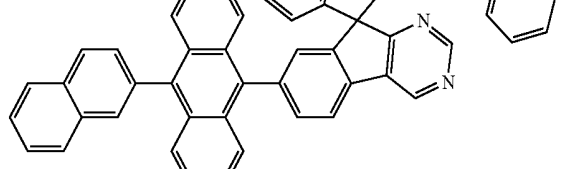
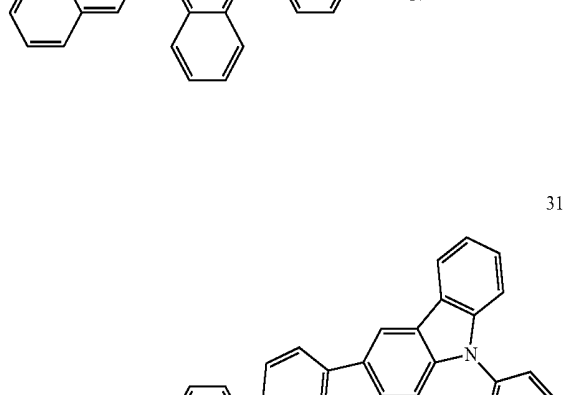
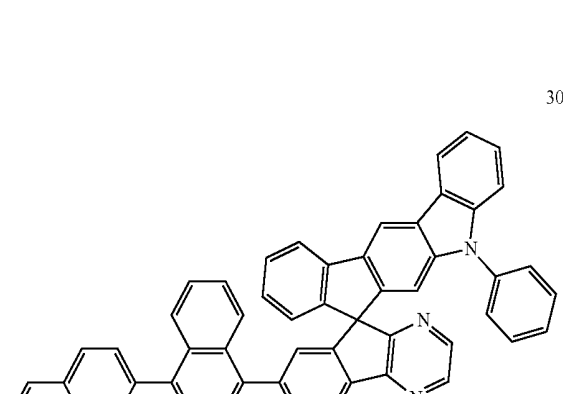
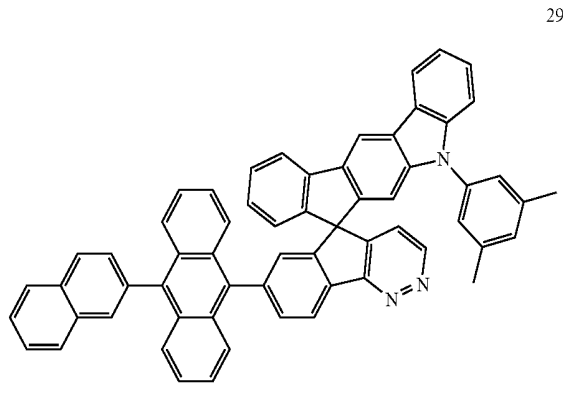
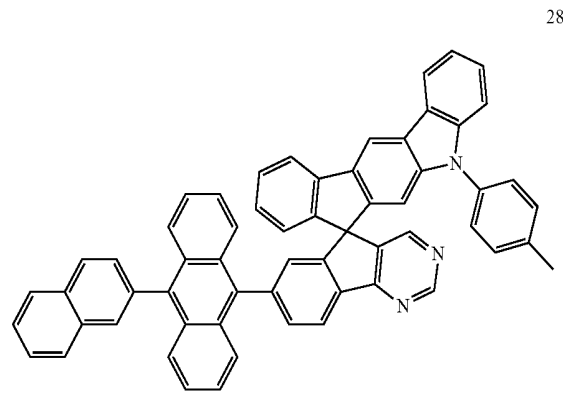
65



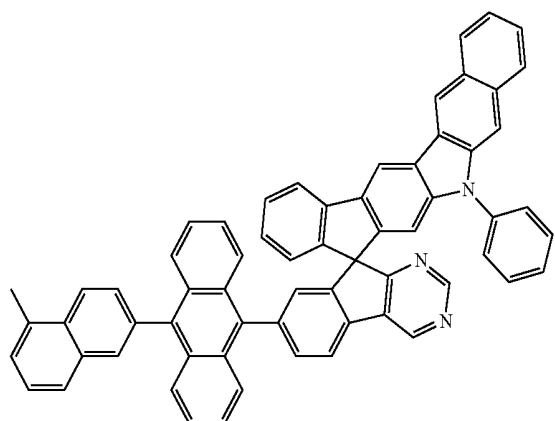
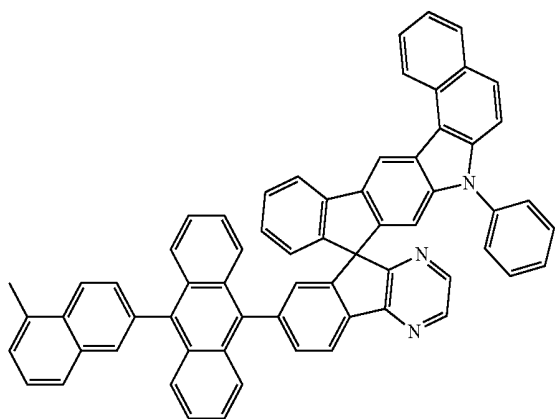
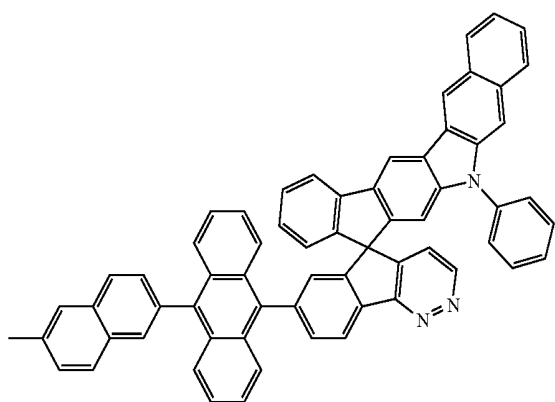
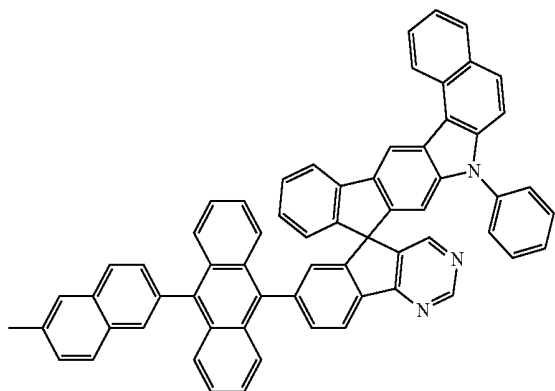
77
-continued



78
-continued



-continued



32 The condensed-cyclic compound represented by Formula 1 has a condensed ring core based on spirobifluorene. Thus, the molecular structure of the condensed-cyclic compound may have strong resistance against electrons, and accordingly, in an organic light-emitting device, deterioration of the molecular structure by electrons may be reduced, which may result in an increase in lifespan of the organic light-emitting device. In addition, the condensed-cyclic compound may have a high triplet (T1) energy level, and accordingly, in an emission layer, triplet excitons may be more prone to collide with one another. Due to triplet-triplet annihilation (TTA) effect, the organic light-emitting device may have improved efficiency.

15 In the condensed-cyclic compound represented by Formula 1, ring A₂ may be selected from a pyridine group, a pyrimidine group, a pyrazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinoxaline group, a quinoxaline group, and a cinnoline group. That is, the condensed-cyclic compound may have at least one nitrogen atom in the condensed ring core based on spirobifluorene. Therefore, the condensed-cyclic compound represented by Formula 1 may have improved electron injecting characteristics into the emission layer, and thus formation of excitons may be more prone to occur in the emission layer. Consequently, efficiency of the organic light-emitting device may be improved.

33 In the condensed-cyclic compound represented by Formula 1, L₁ and L₂ may each independently be selected from substituted or unsubstituted condensed polycyclic groups not including heteroatoms as ring-forming atoms, in which 3 or more carbocyclic groups are condensed with each other. Both a₁ and a₂ are not 0, wherein a₁ and a₂ each denote the number of L₁ (s) and L₂ (s), respectively. That is, in Formula 1, L₁ may be essentially present in a group represented by 34 *-(L₁)_{a1}-(R₁)_{b1}], and L₂ may be essentially present in a group represented by *-(L₂)_{a2}-(R₂)_{b2}]. In Formula 1, a sum of c₁ and c₂ is 1 or greater. That is, at least one selected from a group represented by *-(L₁)_{a1}-(R₁)_{b1}] and a group represented by *-(L₂)_{a2}-(R₂)_{b2}] may be essentially present in 35 Formula 1. Therefore, when the condensed-cyclic compound represented by Formula 1 is used, a suitable energy level between a host and a dopant may be obtained, and exciton energy generated from the host may be efficiently transferred, and thus the organic light-emitting device may 45 have improved efficiency.

The condensed cyclic compound represented by Formula 1 may be synthesized by using a known organic synthetic method. A method of synthesizing the condensed cyclic compound may be recognizable by one of ordinary skill in 46 the art in view of the following embodiments.

50 The at least one condensed cyclic compound of Formula 1 may be used between a pair of electrodes of an organic light-emitting device. For example, the condensed cyclic compound may be included in an emission layer. In one or more embodiments, the condensed cyclic compound of Formula 1 may be used as a material for a forming capping layer positioned outside a pair of electrodes of an organic light-emitting device.

60 According to an aspect, an organic light-emitting device includes: a first electrode, a second electrode, and an organic layer between the first electrode and the second electrode, wherein the organic layer may include an emission layer and at least one condensed cyclic compound represented by Formula 1 described above.

65 The expression “(an organic layer) includes at least one condensed-cyclic compound” as used herein refer to an embodiment in which “(an organic layer) includes identical

compounds represented by Formula 1 and to an embodiment in which (an organic layer) includes two or more different condensed-cyclic compounds represented by Formula 1.”

In some embodiments, the organic layer may include only Compound 1 as the condensed-cyclic compound. In this case, Compound 1 may be included in an emission layer of the organic light-emitting device. In one or more embodiments, the organic layer may include both Compounds 1 and 2 as the condensed-cyclic compound. In this regard, Compounds 1 and 2 may be situated in the same layer (for example, Compounds 1 and 2 may all be situated in an emission layer), or in different layers (for example, Compound 1 may be situated in a hole transport layer, and Compound 2 may be situated in an emission layer).

The organic layer may include i) a hole transport region disposed between the first electrode (anode) and the emission layer, the hole transport region including at least one of a hole injection layer, a hole transport layer, an emission auxiliary layer, and an electron blocking layer, and ii) an electron transport region disposed between the emission layer and the second electrode (cathode), the electron transport region including at least one of a hole blocking layer, an electron transport layer, and an electron injection layer. At least one of the hole transport region and the emission layer may include the at least one condensed cyclic compound represented by Formula 1. For example, the emission layer of the organic light-emitting device may include the at least one condensed cyclic compound represented by Formula 1. The at least one condensed cyclic compound represented by Formula 1 included in the emission layer may act as a host, and the emission layer may further include a dopant. The dopant may be a phosphorous dopant or a fluorescent dopant. In an embodiment, the dopant may be a fluorescent dopant.

The organic light-emitting device may further include at least one of a first capping layer disposed on a pathway through which light extracted or emitted from an emission layer proceeds toward the outside through the first electrode; and a second capping layer disposed on a pathway through which light extracted or emitted from an emission layer proceeds toward the outside through the second electrode, wherein the at least one of the first capping layer and the second capping layer may include the at least one condensed-cyclic compound.

In some embodiments, an organic light-emitting device may have i) a stack structure including a first electrode, an organic layer, a second electrode, and a second capping layer, which are sequentially stacked in this stated order, ii) a stack structure including a first capping layer, a first electrode, an organic layer, and a second electrode, which are sequentially stacked in this stated order, or iii) a stack structure including a first capping layer, a first electrode, an organic layer, a second electrode, and a second capping layer, which are sequentially stacked in this stated order, wherein at least one of the first capping layer and the second capping layer may include the at least one condensed-cyclic compound.

The term “organic layer” as used herein refers to a single layer and/or a plurality of layers disposed between the first electrode and the second electrode of the organic light-emitting device. Materials included in the “organic layer” are not limited to organic materials.

Description of FIG. 1

FIG. 1 is a schematic cross-sectional 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.

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

First Electrode 110

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

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 the first electrode 110 may be selected from materials with a high work function to facilitate hole injection.

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 the first electrode 110 may be selected from indium tin oxide (ITO), indium zinc oxide (IZO), tin oxide (SnO₂), zinc oxide (ZnO), and any combinations thereof, but embodiments are not limited thereto. In one or more embodiments, when the first electrode 110 is a semi-transmissive electrode or a reflective electrode, the material for the first electrode 110 may be selected from magnesium (Mg), aluminum (Al), aluminum-lithium (Al—Li), calcium (Ca), magnesium-indium (Mg—In), magnesium-silver (Mg—Ag), and any combinations thereof, but embodiments are not limited thereto.

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 ITO/Ag/ITO structure, but embodiments of the structure of the first electrode 110 are not limited thereto.

Organic Layer 150

The organic layer 150 may be disposed on the first electrode 110. The organic layer 150 may include an emission region.

The organic layer 150 may include a hole transport region between the first electrode 110 and the emission region and an electron transport region between the emission region and the second electrode 190.

Hole Transport Region in Organic Layer 150

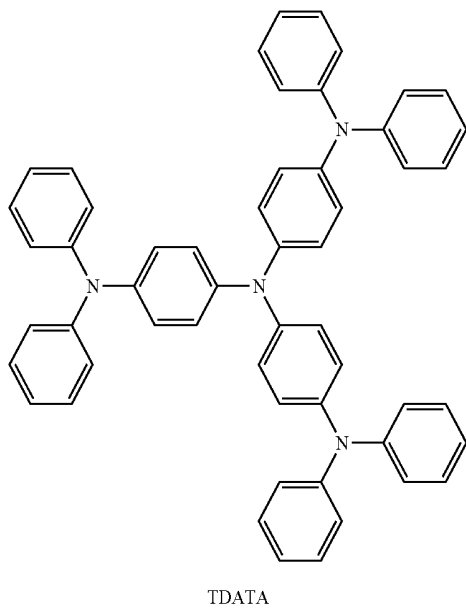
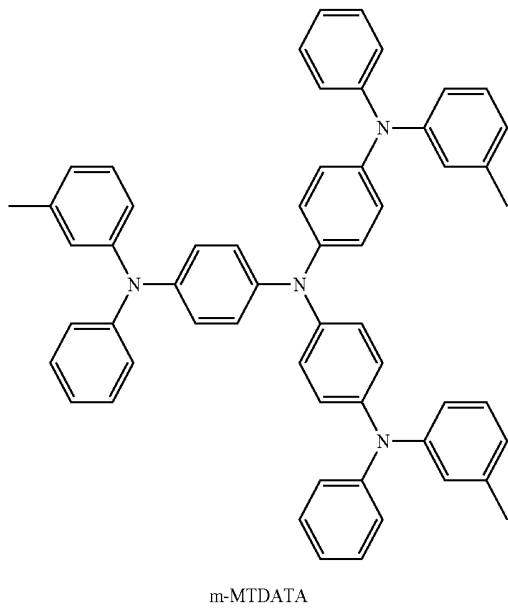
The hole transport region may have i) a single-layered structure including a single layer including a single material, ii) a single-layered structure including a single layer including a plurality of different materials, or iii) a multi-layered structure having a plurality of layers including a plurality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

The hole transport region may include at least one layer selected from a hole injection layer, a hole transport layer, an emission auxiliary layer, and an electron blocking layer.

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 layers of each structure are sequentially stacked on the first electrode 110 in each stated order, but embodiments of the structure of the hole transport region are not limited thereto.

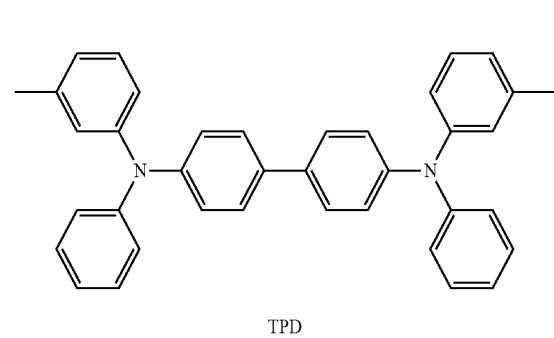
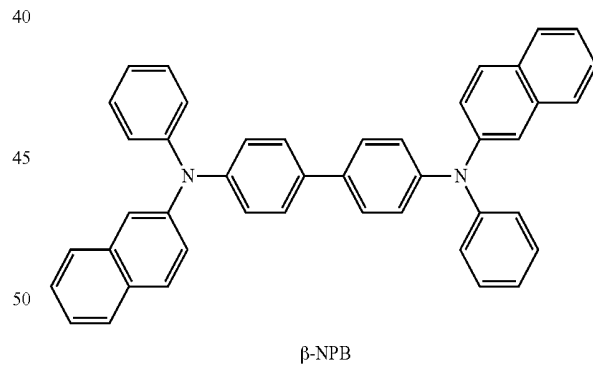
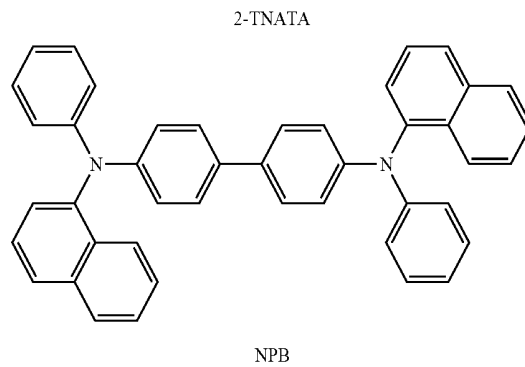
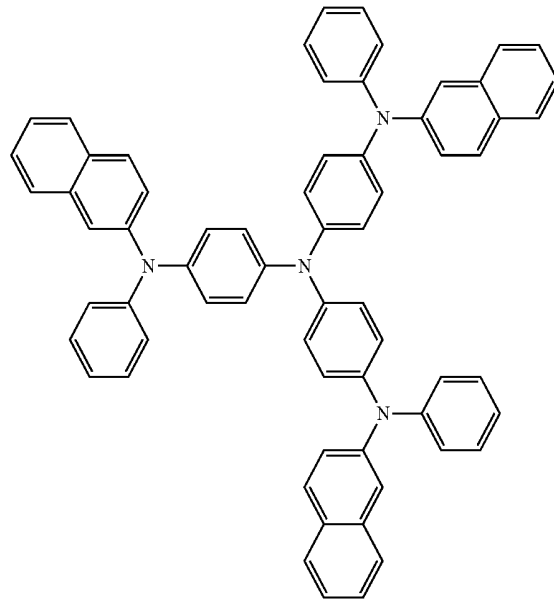
83

The hole transport region may include at least one selected from m-MTDATA, TDATA, 2-TNATA, NPB (NPB), p-NPB, TPD, Spiro-TPD, Spiro-NPB, methylated-NPB, TAPC, HMTPD, 4,4',4''-tris(N-carbazolyl)triphenylamine (TCTA), polyaniline/dodecylbenzenesulfonic acid (PANI/DBSA), poly(3,4-ethylenedioxythiophene)/poly(4-styrenesulfonate) (PEDOT/PSS), polyaniline/camphor sulfonic acid (PANI/CSA), polyaniline/poly(4-styrenesulfonate) (PANI/PSS), a compound represented by Formula 201, and a compound represented by Formula 202, but embodiments are not limited thereto:



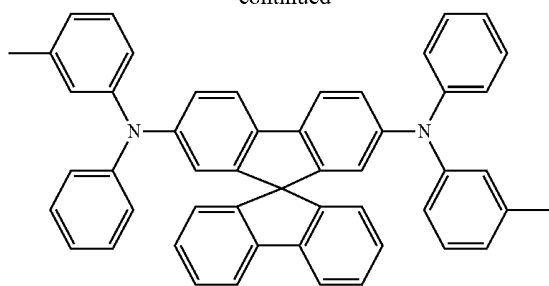
84

-continued

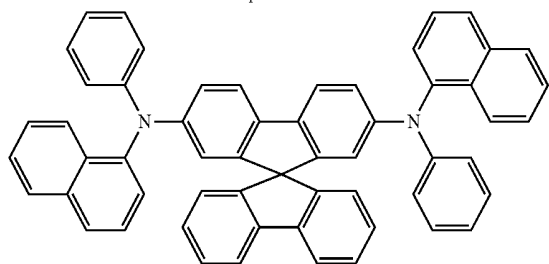


85

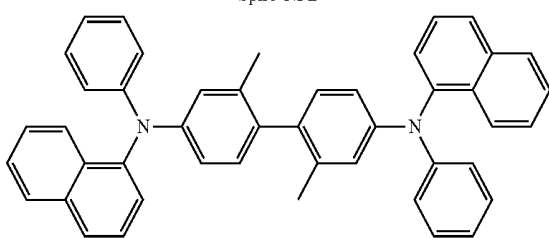
-continued



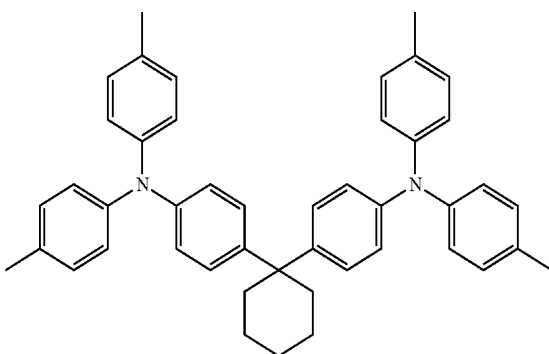
Spiro-TPD



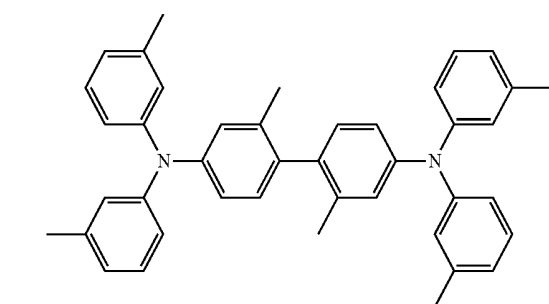
Spiro-NPB



methylated NPB



TAPC



HMTPD

86

-continued

5

10

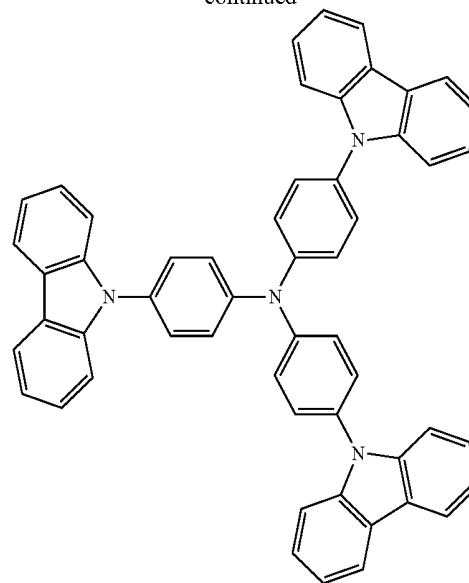
15

20

25

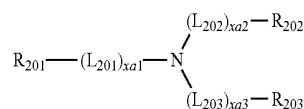
30

35

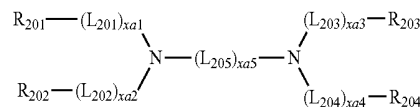


TCTA

Formula 201



Formula 202



wherein, in Formulae 201 and 202,

L₂₀₁ to L₂₀₄ may each independently be selected from a substituted or unsubstituted C₃-C₁₀ cycloalkylene group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkylene group, a substituted or unsubstituted C₃-C₁₀ cycloalkenylene group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkenylene group, a substituted or unsubstituted C₆-C₆₀ arylene group, a substituted or unsubstituted C₁-C₆₀ heteroarylene group, a substituted or unsubstituted divalent non-aromatic condensed polycyclic group, and a substituted or unsubstituted divalent non-aromatic condensed heteropolycyclic group,

L₂₀₅ may be selected from *-O-*, *-S-*, *-N(Q₂₀₁)-*, a substituted or unsubstituted C₁-C₂₀ alkylene group, a substituted or unsubstituted C₂-C₂₀ alkenylene group, a substituted or unsubstituted C₃-C₁₀ cycloalkylene group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkylene group, a substituted or unsubstituted C₃-C₁₀ cycloalkenylene group, a substituted or unsubstituted C₁-C₁₀ heterocycloalkenylene group, a substituted or unsubstituted C₆-C₆₀ arylene group, a substituted or unsubstituted C₁-C₆₀ heteroarylene group, a substituted or unsubstituted divalent non-aromatic condensed polycyclic group, and a substituted or unsubstituted divalent non-aromatic condensed heteropolycyclic group,

x_{a1} to x_{a4} may each independently be an integer from 0 to 3,

x_{a5} may be an integer from 1 to 10, and

R₂₀₁ to R₂₀₄ and Q₂₀₁ may each independently be selected from 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, and a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group.

In some embodiments, in Formula 202, R₂₀₁ and R₂₀₂ may optionally be bound via a single bond, a dimethyl-methylene group, or a diphenyl-methylene group, and R₂₀₃ and R₂₀₄ may optionally be bound via a single bond, a dimethyl-methylene group, or a diphenyl-methylene group.

In one embodiment, in Formulae 201 and 202, L₂₀₁ to L₂₀₅ may each independently be selected from

a phenylene group, a pentalenylene group, an indenylene group, a naphthylene group, an azulenylenylene group, a heptalenylene group, an indacenylenylene group, an acenaphthylene group, a fluorenylenylene group, a spiro-bifluorenylenylene group, a benzofluorenylenylene group, a dibenzofluorenylenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylenylene group, a fluoranthenylenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a naphthacenylenylene group, a picenylene group, a perylenylene group, a pentaphenylenylene group, a hexacenylenylene group, a pentacenylenylene group, a rubicenylenylene group, a coronenylenylene group, an ovalenylenylene 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, and a pyridinylenylene group; and

a phenylene group, a pentalenylene group, an indenylene group, a naphthylene group, an azulenylenylene group, a heptalenylene group, an indacenylenylene group, an acenaphthylene group, a fluorenylenylene group, a spiro-bifluorenylenylene group, a benzofluorenylenylene group, a dibenzofluorenylenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylenylene group, a fluoranthenylenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a naphthacenylenylene group, a picenylene group, a perylenylene group, a pentaphenylenylene group, a hexacenylenylene group, a pentacenylenylene group, a rubicenylenylene group, a coronenylenylene group, an ovalenylenylene 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, and a pyridinylenylene 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 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 C₁-C₁₀ alkyl group, a phenyl group substituted with —F, a pentalenyl group, an indenyl group, a naphthyl group, a fluorenyl 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, 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 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 C₁-C₁₀ alkyl group, a phenyl group substituted with —F, a pentalenyl group, an indenyl group, a naphthyl group, a fluorenyl 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

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, a pyridinyl group, —Si(Q₃₁)(Q₃₂)(Q₃₃), and —N(Q₃₁)(Q₃₂).

wherein 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.

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

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

In one or more embodiments, R₂₀₁ to R₂₀₄ and Q₂₀₁ may each independently be selected from a phenyl group, a biphenyl group, a terphenyl group, a pentalenyl group, an indenyl group, a naphthyl 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; and

a phenyl group, a biphenyl group, a terphenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenylenylene group, a heptalenyl group, an indacenylenylene 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, 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 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 C₁-C₁₀ alkyl group, a phenyl group substituted with —F, a pentalenyl group, an indenyl group, a naphthyl group, a fluorenyl 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, a pyridinyl group, $-\text{Si}(\text{Q}_{31})(\text{Q}_{32})(\text{Q}_{33})$, and $-\text{N}(\text{Q}_{31})(\text{Q}_{32})$,

wherein Q_{31} to Q_{33} may be the same as those described herein.

In one or more embodiments, in Formula 201, at least one of R_{201} to R_{203} may be selected from

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

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, but embodiments are not limited thereto.

In one or more embodiments, in Formula 202, i) R_{201} and R_{202} may be bound via a single bond, and/or ii) R_{203} and R_{204} may be bound via a single bond.

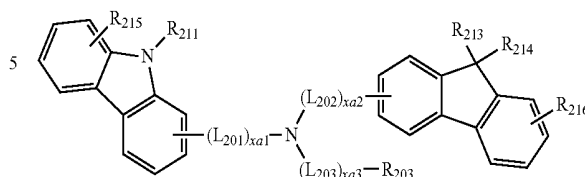
In one or more embodiments, in Formula 202, at least one of R_{201} to R_{204} may be selected from

a carbazolyl group; and

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, but embodiments are not limited thereto.

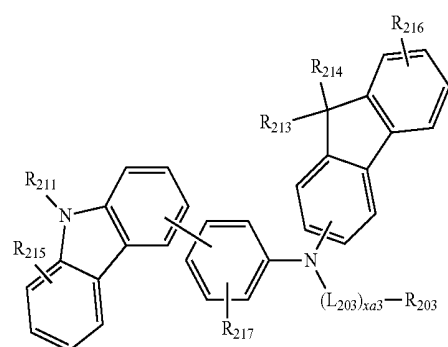
The compound represented by Formula 201 may be represented by Formula 201A:

Formula 201A



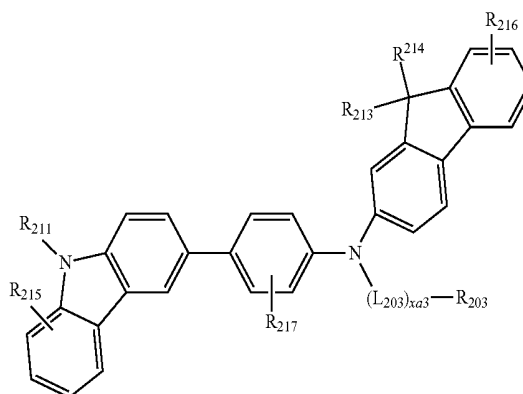
In some embodiments, the compound represented by Formula 201 may be represented by Formula 201A(1), but embodiments are not limited thereto:

Formula 201A(1)



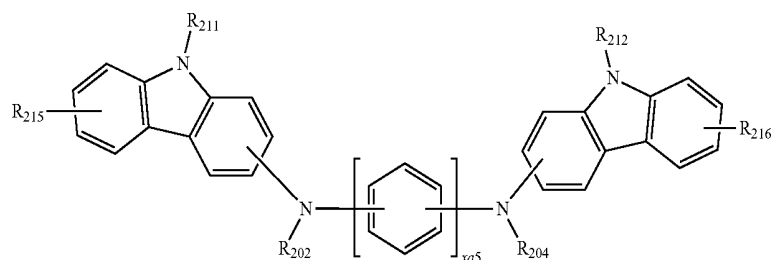
In some embodiments, the compound represented by Formula 201 may be represented by Formula 201A-1, but embodiments are not limited thereto:

Formula 201A-1



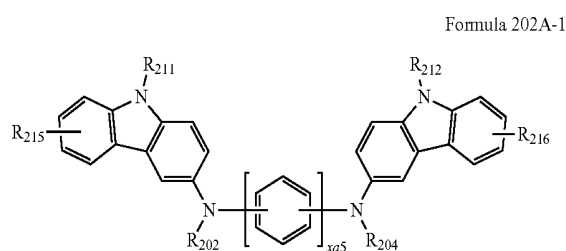
In some embodiments, the compound represented by Formula 202 may be represented by Formula 202A:

Formula 202A



91

In some embodiments, the compound represented by Formula 202 may be represented by Formula 202A-1:



In Formulae 201A, 201A(1), 201A-1, 202A, and 202A-1, L₂₀₁ to L₂₀₃, xa1 to xa3, and R₂₀₂ to R₂₀₄ may be the same as those described herein,

R₂₁₁ and R₂₁₂ may each be the same as described above with reference to R₂₀₃, and

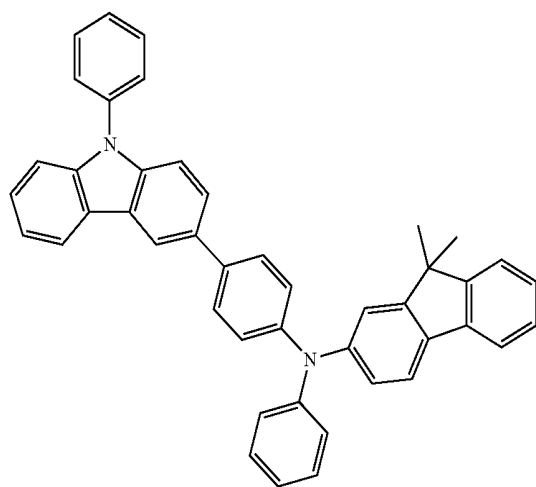
R₂₁₃ to R₂₁₇ may each independently be selected from hydrogen, deuterium, —F, —Cl, —Br, —I, a hydroxyl

92

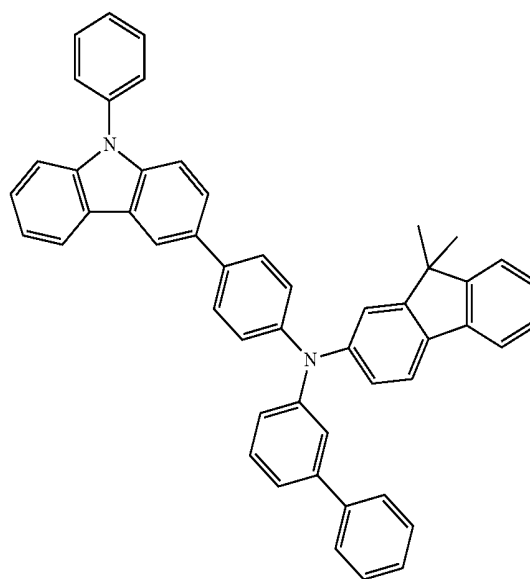
group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C₁-C₂₀ alkyl group, a C₁-C₂₀ alkoxy group, a cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl group, a biphenyl group, a terphenyl group, a phenyl group substituted with a C₁-C₁₀ 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.

The hole transport region may include at least one compound selected from Compounds HT1 to HT39, but embodiments are not limited thereto:

HT1



HT2



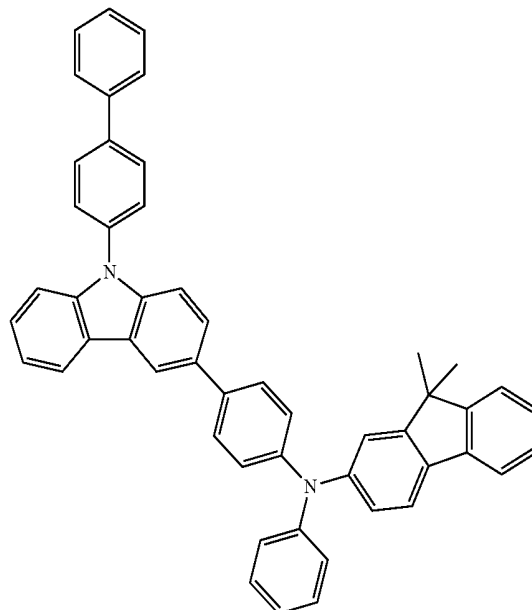
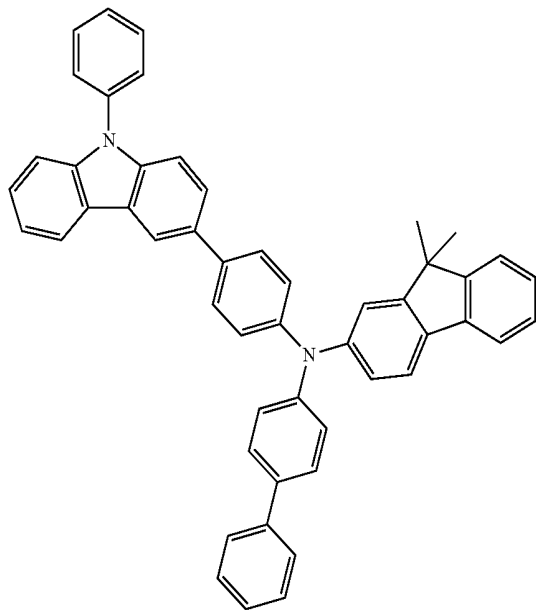
93

94

-continued

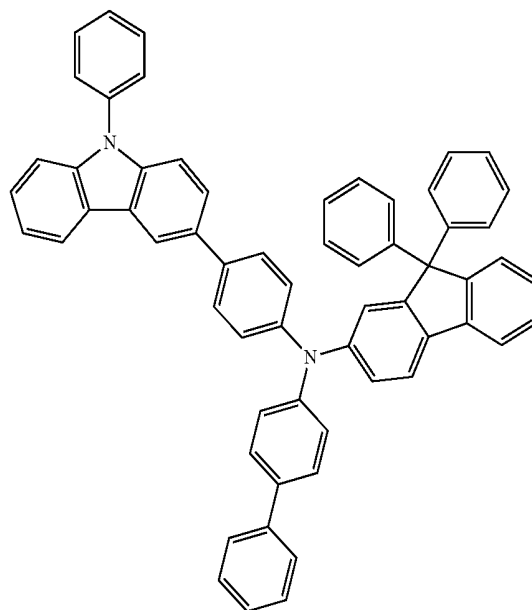
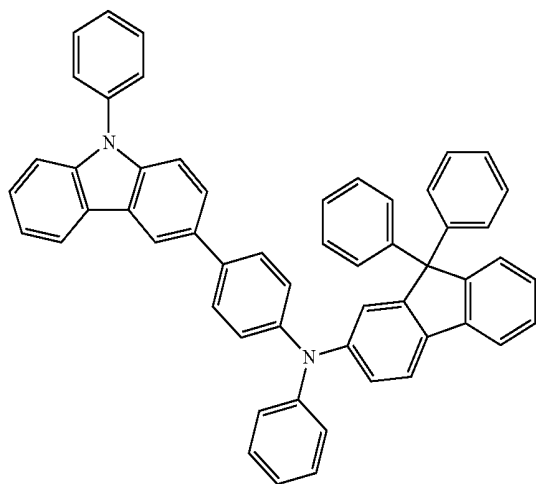
HT3

HT4



HT5

HT6



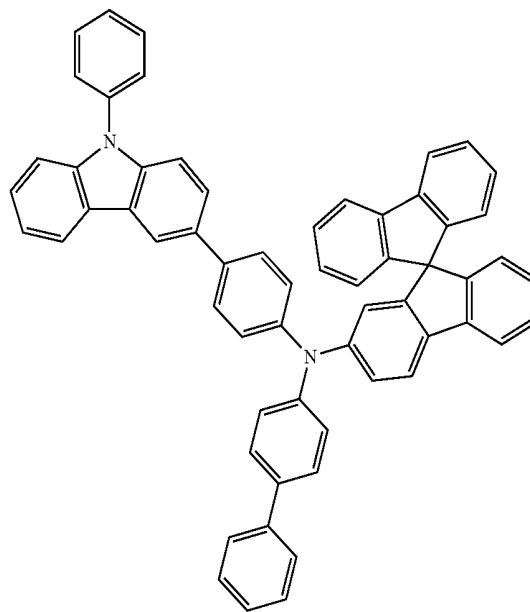
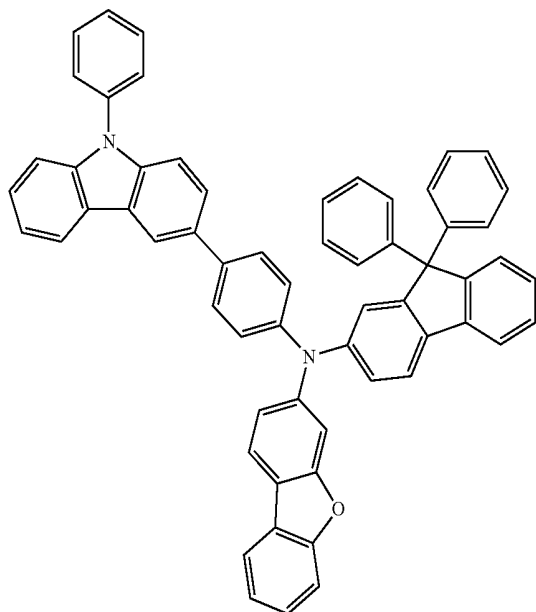
95

96

-continued

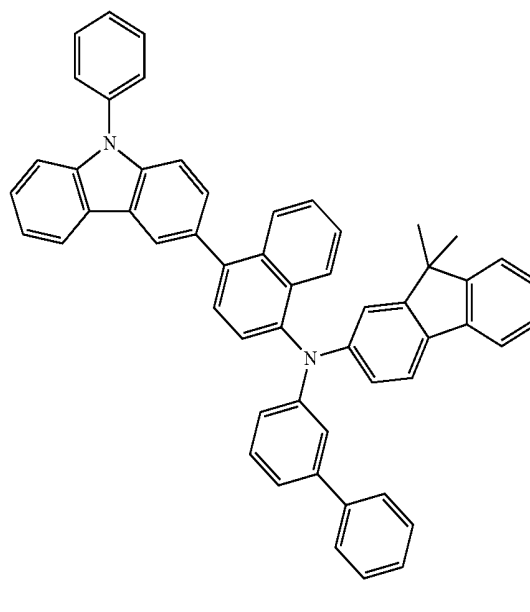
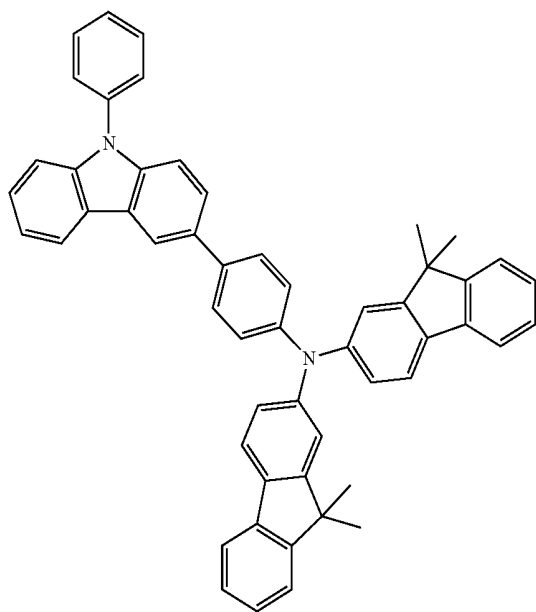
HT7

HT8



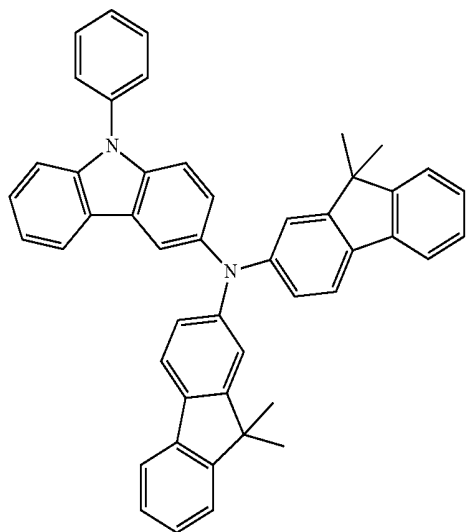
HT-9

HT-10



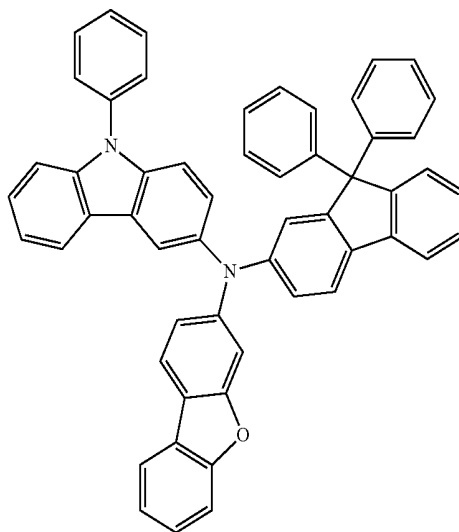
97

-continued
HT11

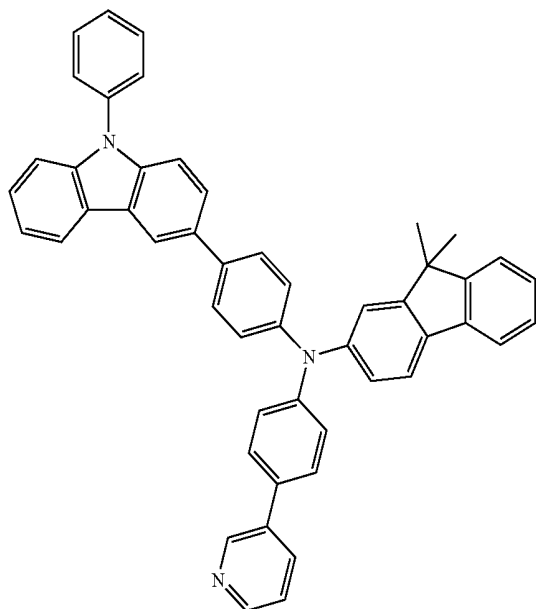


98

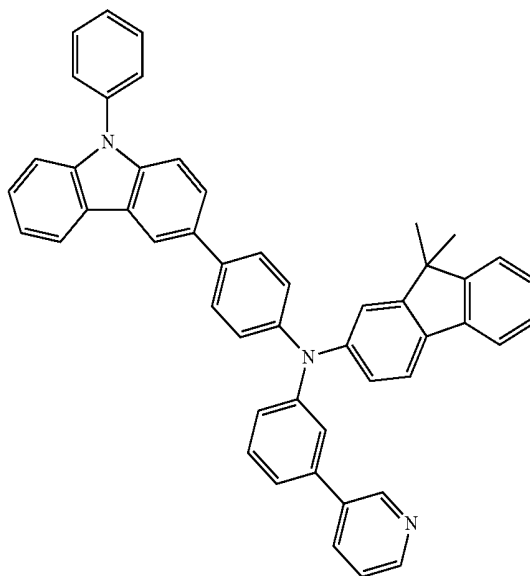
HT12



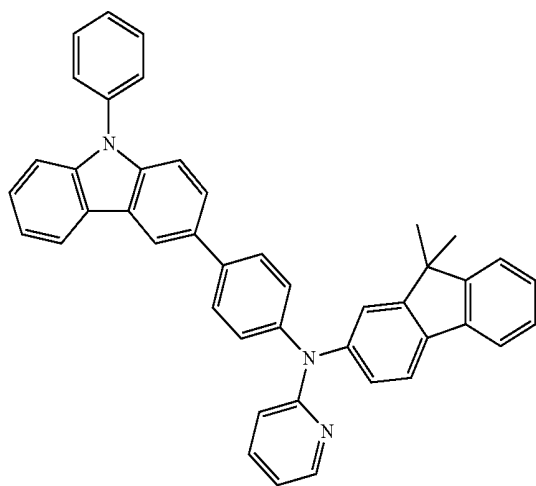
HT13



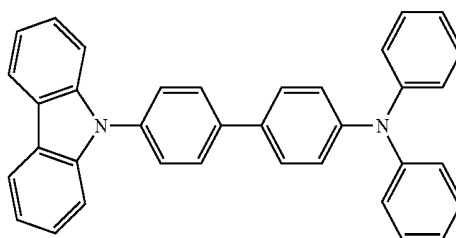
HT14



HT15



HT16



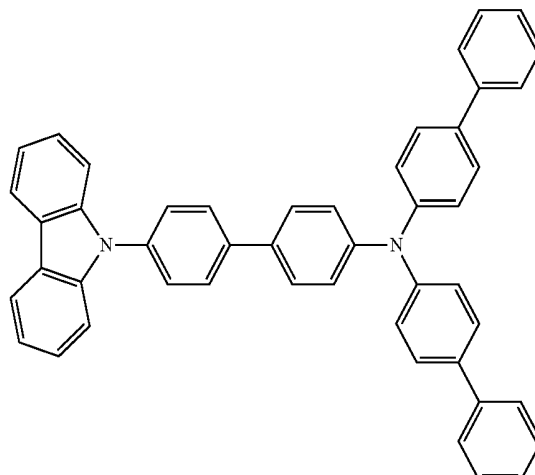
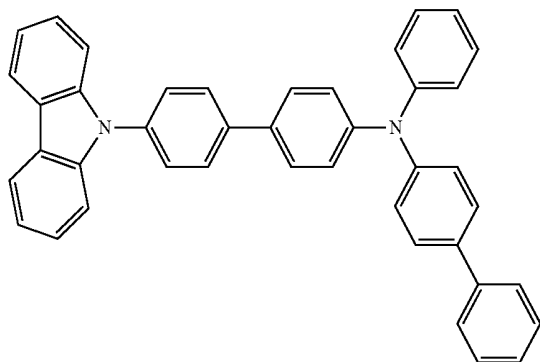
99

100

-continued

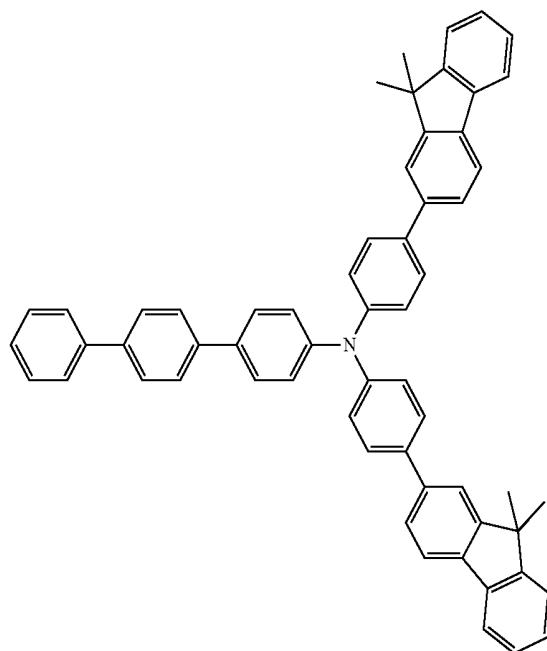
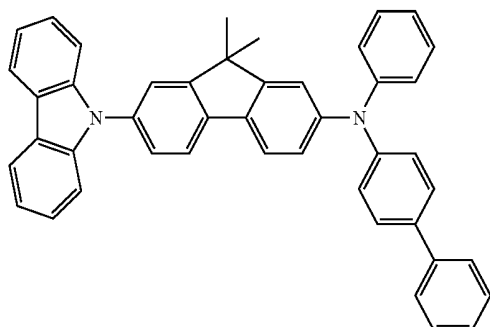
HT17

HT18



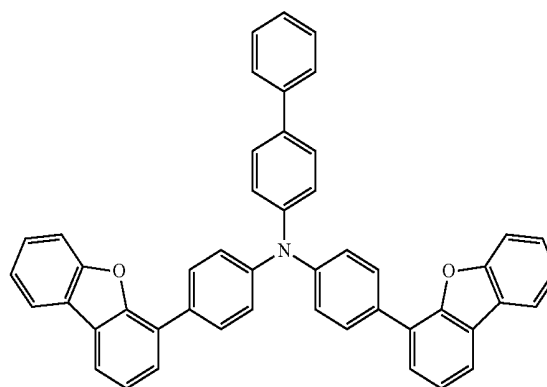
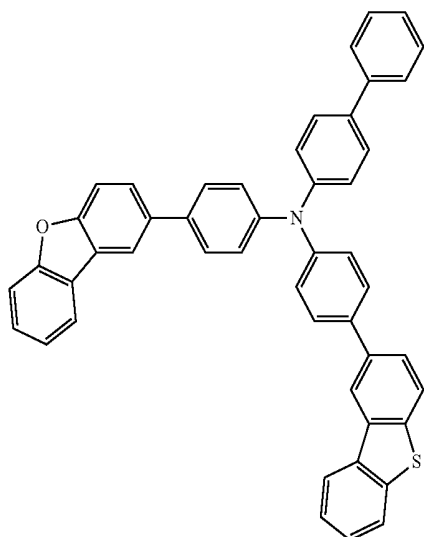
HT19

HT20



HT21

HT22

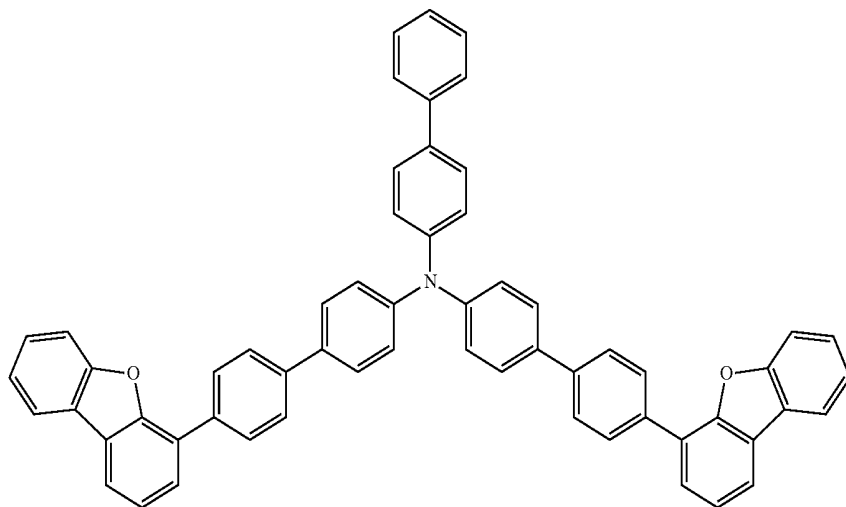


101

102

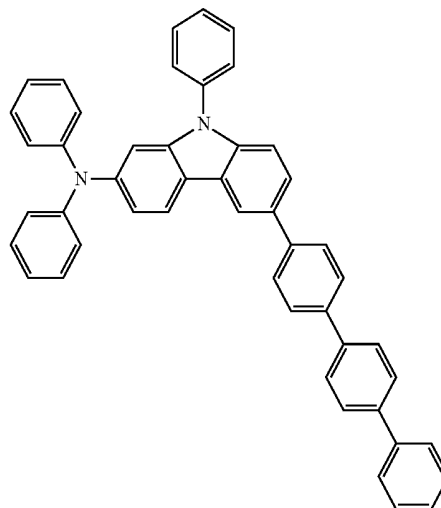
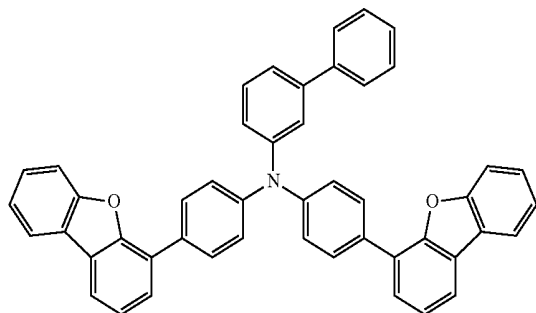
-continued

HT23



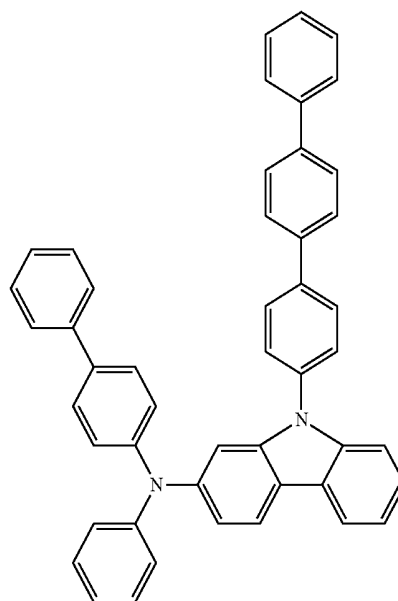
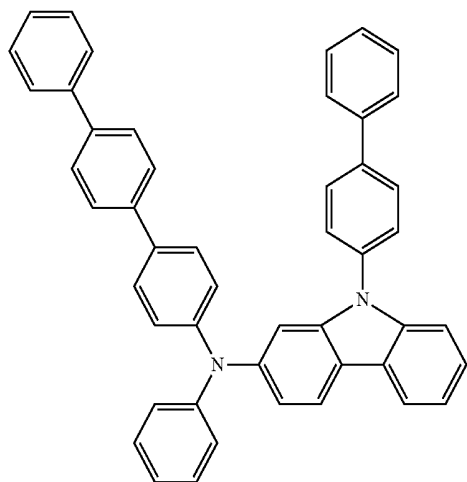
HT24

HT25



HT26

HT27



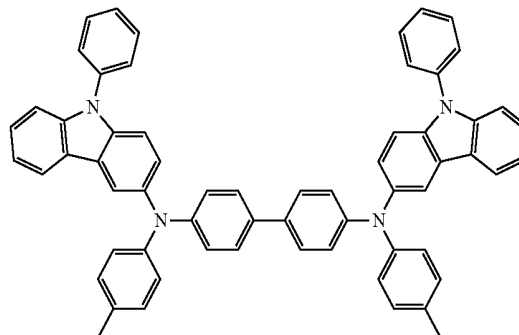
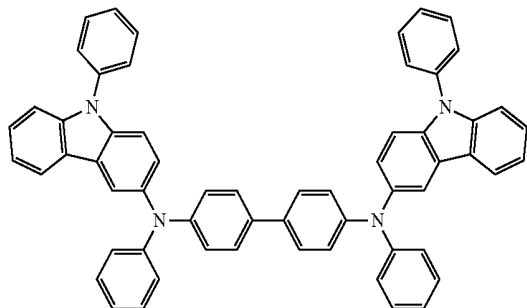
103

104

-continued

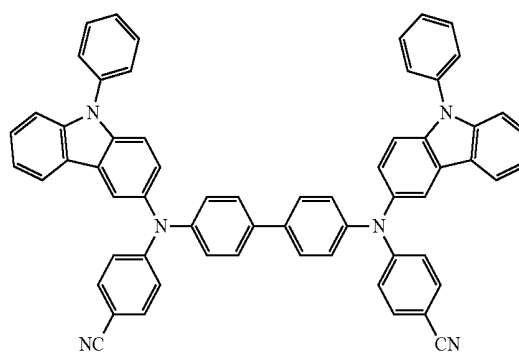
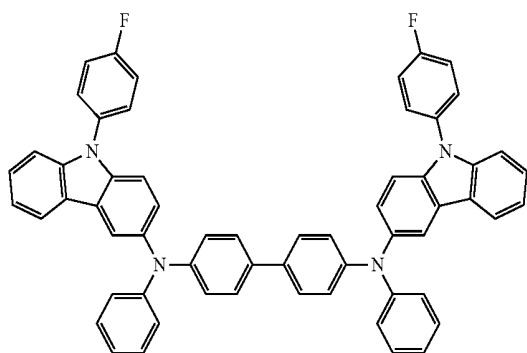
HT28

HT29



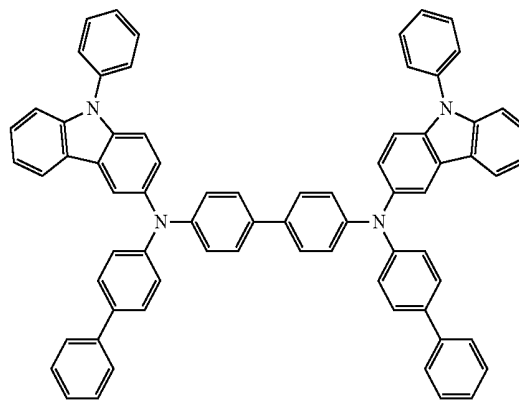
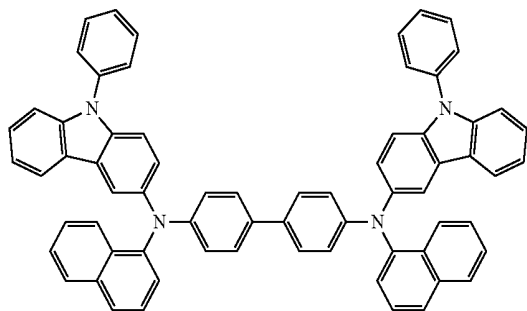
HT30

HT31



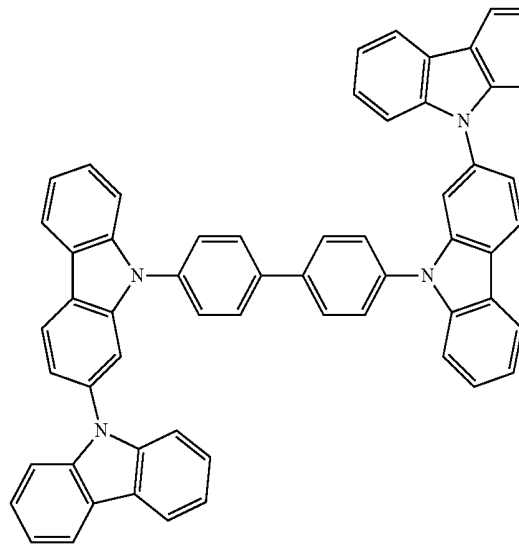
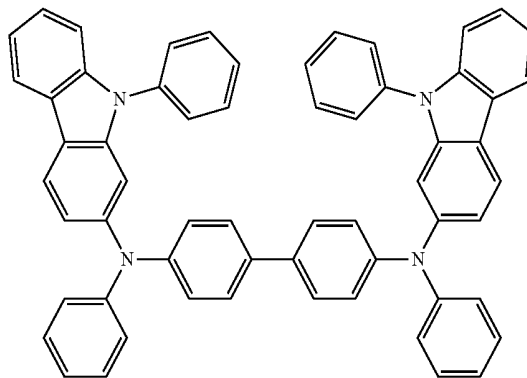
HT32

HT33



HT34

HT35

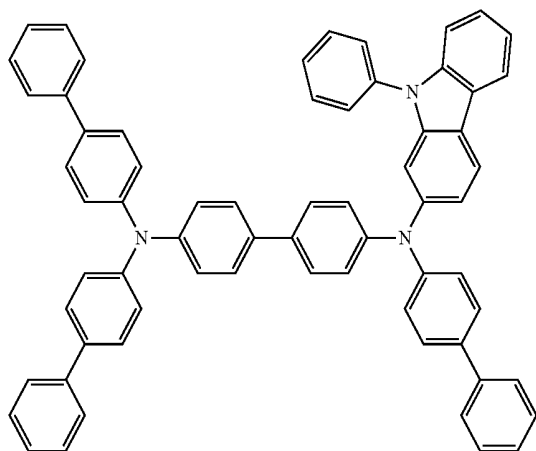


105

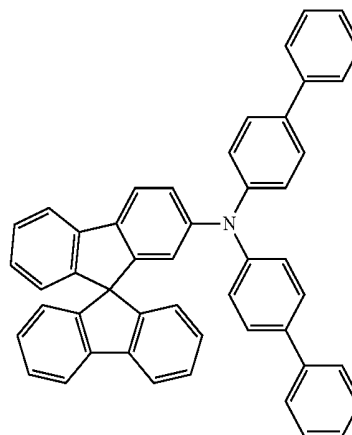
106

-continued
HT36

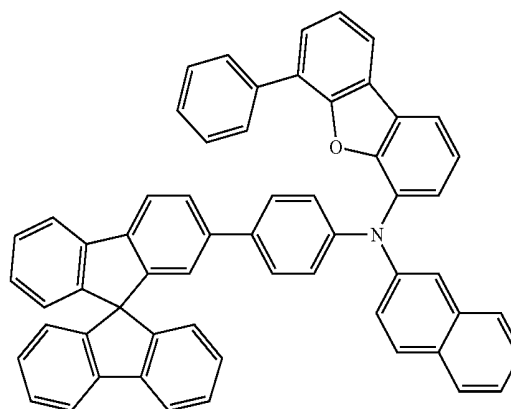
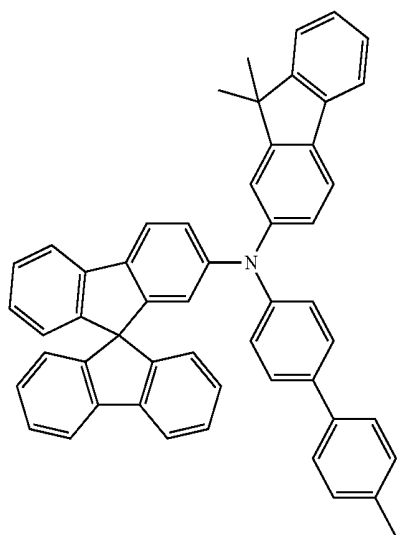
HT37



HT38



HT39



The 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, 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, 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 any of these ranges, satisfactory hole transporting characteristics may be obtained without a substantial increase in driving voltage.

The emission auxiliary layer may increase the light-emission efficiency by compensating for an optical resonance distance depending on the wavelength of light emitted by an emission layer, and the electron blocking layer may block or reduce 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

The hole transport region may further include, in addition to these materials, a charge generating material to improve

conductive properties. The charge generating material may be homogeneously or non-homogeneously dispersed in the hole transport region.

The charge generating material may be, for example, a p-dopant.

In one embodiment, the p-dopant may have a lowest unoccupied molecular orbital (LUMO) level of about -3.5 eV or less.

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

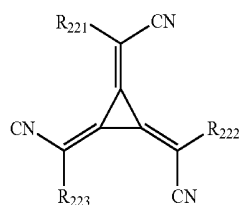
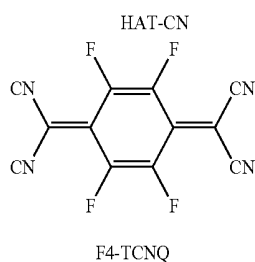
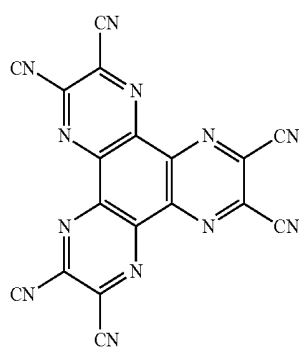
In some embodiments, the p-dopant may include at least one selected from

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

a metal oxide, such as tungsten oxide or molybdenum oxide;

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

a compound represented by Formula 221, but embodiments of the present disclosure are not limited thereto:



Formula 221

wherein, in Formula 221,

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, wherein at least one selected from R_{221} to R_{223} may include 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

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, wherein 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.

In one embodiment, the emission layer of the organic light-emitting device **10** may be a first color light emission layer,

the organic light-emitting device **10** may further include i) at least one second color light emission layer or ii) at least one second color light emission layer and at least one third color light emission layer, between the first electrode **110** and the second electrode **190**,

a maximum emission wavelength of the first color light emission layer, a maximum emission wavelength of the second color light emission layer, and a maximum emission wavelength of the third color light emission layer are identical to or different from each other, and

the organic light-emitting device **10** may emit mixed light including the first color light and the second color light, or mixed light including the first color light, the second color light, and the third color light, but embodiments are not limited thereto.

For example, the maximum emission wavelength of the first color light emission layer is different from a maximum emission wavelength of the second color light emission layer, and the mixed light including first color light and second color light may be white light, but embodiments are not limited thereto.

In one or more embodiments, the maximum emission wavelength of the first color light emission layer, the maximum emission wavelength of the second color light emission layer, and the maximum emission wavelength of the third color light emission layer may be different from one another, and the mixed light including the first color light, the second color light, and the third color light may be white light. However, embodiments are not limited thereto.

The emission layer may include a host and a dopant. The dopant may include a phosphorous dopant or a fluorescent dopant.

The amount of the dopant in the emission layer may be in a range of about 0.01 parts to about 15 parts by weight based on 100 parts by weight of the host, but embodiments are not limited thereto.

The thickness of the emission layer may be in a range of about 100 Å to about 1,000 Å, and in some embodiments, about 200 Å to about 600 Å. When the thickness of the emission layer is within any of these ranges, excellent light-emission characteristics may be obtained without a substantial increase in driving voltage.

Host in Emission Layer

The host may include the condensed cyclic compound represented by Formula 1.

In one or more embodiments, the host may include a compound represented by Formula 301:



wherein, in Formula 301,

Ar_{301} may be selected from a substituted or unsubstituted C_5 - C_{60} carbocyclic group and a substituted or unsubstituted C_1 - C_{60} heterocyclic group, $xb11$ may be 1, 2, or 3,

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,

$xb1$ may be an integer from 0 to 5,

R_{301} may be selected from deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an

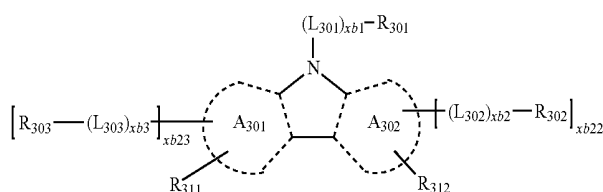
amidino group, a hydrazino group, a hydrazono group, a substituted or unsubstituted C₁-C₆₀ alkyl group, a substituted or unsubstituted C₂-C₆₀ alkenyl group, a substituted or unsubstituted C₂-C₆₀ alkynyl group, a substituted or unsubstituted C₁-C₆₀ alkoxy group, a substituted or unsubstituted C₃-C₁₀ cycloalkyl 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

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),

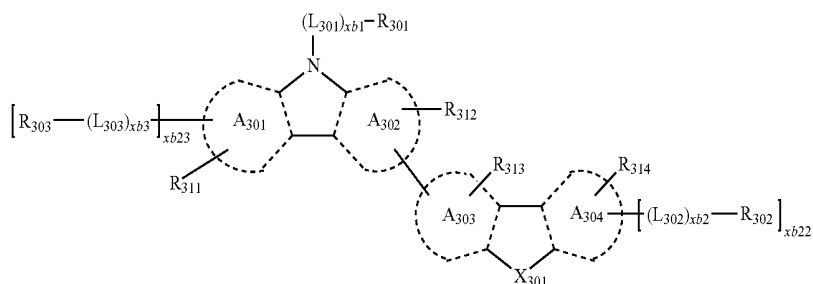
wherein Q₃₁ to Q₃₃ may each independently be selected from a C₁-C₁₀ alkyl group, a C₁-C₁₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group, but embodiments are not limited thereto.

When xb11 in Formula 301 is 2 or greater, at least two Ar₃₀₁ (s) may be bound via a single bond.

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

non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, —Si(Q₃₀₁)(Q₃₀₂)(Q₃₀₃), —N(Q₃₀₁)(Q₃₀₂), —B(Q₃₀₁)(Q₃₀₂), —C(=O)(Q₃₀₁), —S(=O)₂(Q₃₀₁), and —P(=O)(Q₃₀₁)(Q₃₀₂), and

xb21 may be an integer from 1 to 5,

wherein Q₃₀₁ to Q₃₀₃ 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, but embodiments are not limited thereto.

In one embodiment, in Formula 301, Ar₃₀₁ may be selected from

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

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

wherein, in Formulae 301-1 to 301-2,

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,

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

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).

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

L₃₀₁, xb1, R₃₀₁, and Q₃₁ to Q₃₃ may be the same as those described herein,

L₃₀₂ to L₃₀₄ may each independently be substantially the same as described herein with reference to L₃₀₁,

xb2 to xb4 may each independently be substantially the same as described herein with reference to xb1, and

113

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

wherein Q₃₁ to Q₃₃ may be the same as those described above.

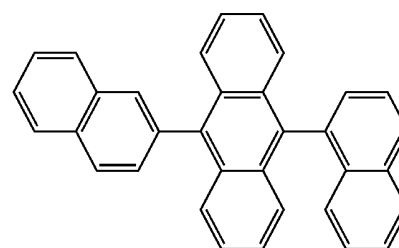
In some embodiments, the host may include an alkaline earth metal complex.

For example, the host may include a beryllium (Be) complex, for example, Compound H55, a magnesium (Mg) complex, or a zinc (Zn) complex.

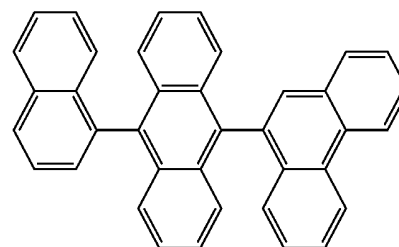
The host may include at least one selected from 9,10-di(2-naphthyl)anthracene (ADN), 2-methyl-9,10-bis(naphthalen-2-yl)anthracene (MADN), 9,10-di-(2-naphthyl)-2-t-butyl-anthracene (TBADN), 4,4'-bis(N-carbazolyl)-1,1'-biphenyl (CBP), 1,3-di-9-carbazolylbenzene (mCP), 1,3,5-tri(carbazol-9-yl)benzene (TCP),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

114

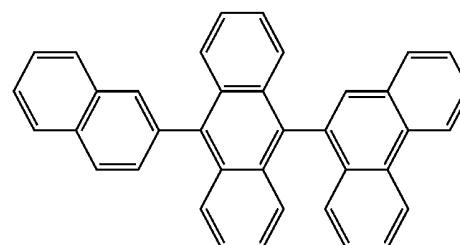
-continued



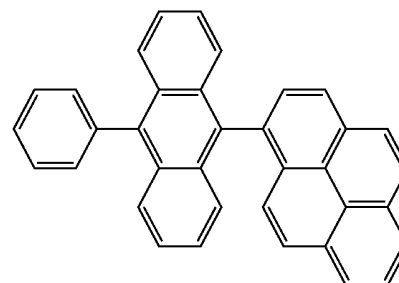
H2



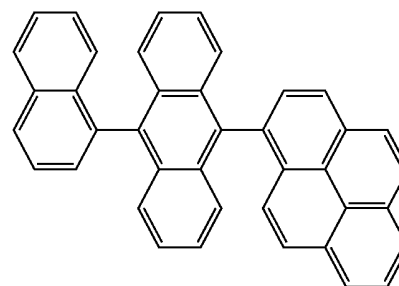
H3



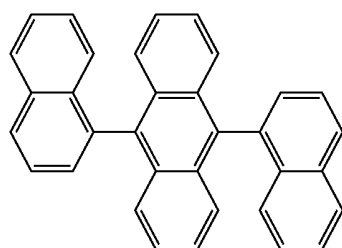
H4



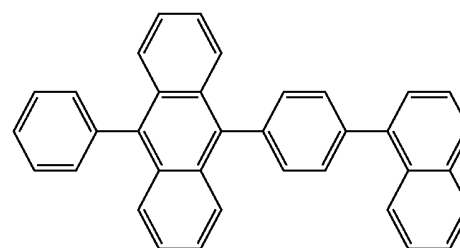
H5



H6



H1

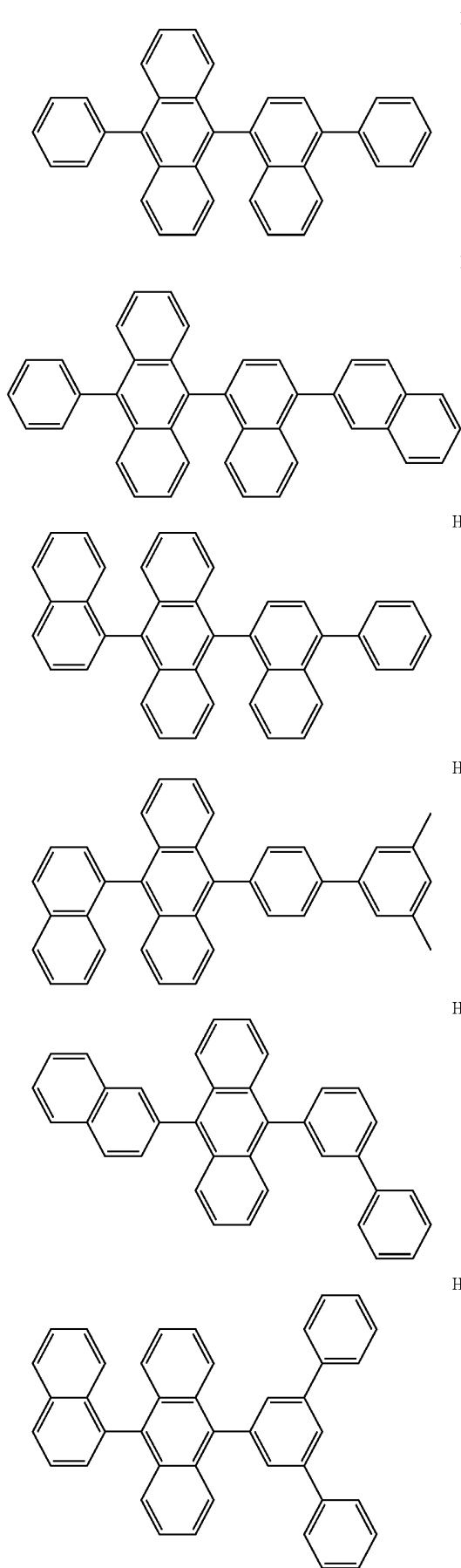


H7

65

115

-continued



H8

5

H9

10

15

20

H10

25

H11

35

40

H12

45

50

H13

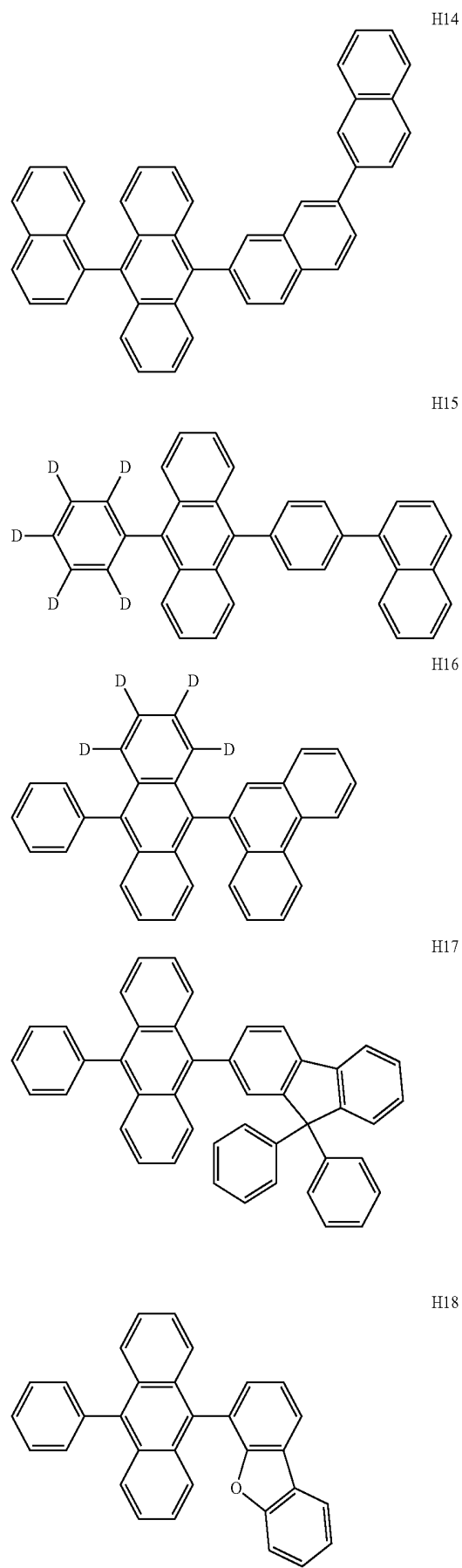
55

60

65

116

-continued



H14

H15

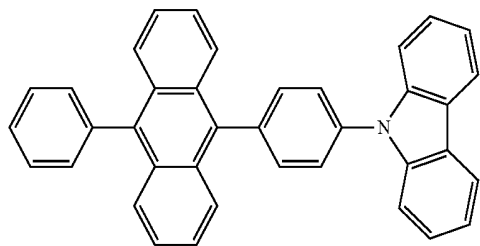
H16

H17

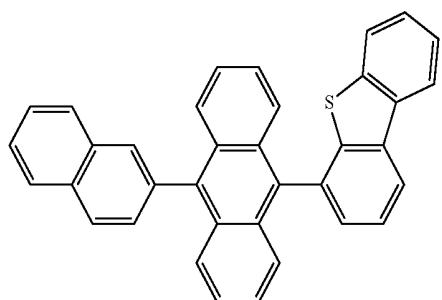
H18

117
-continued

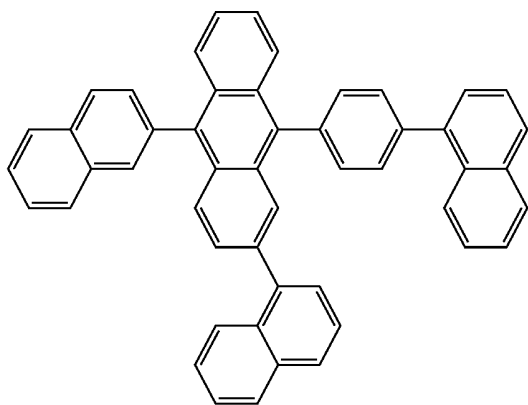
H19



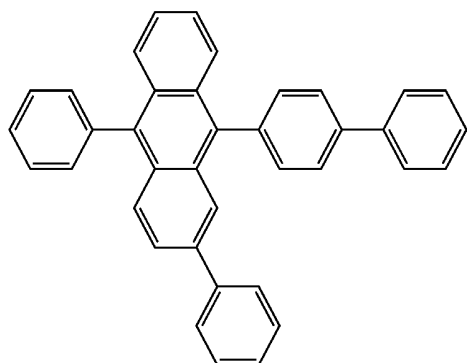
15
HT20



25
HT21

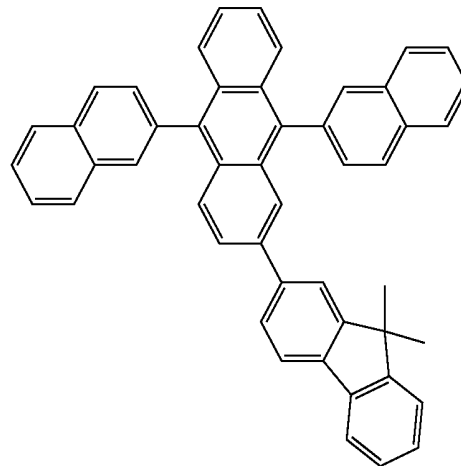


45
HT22

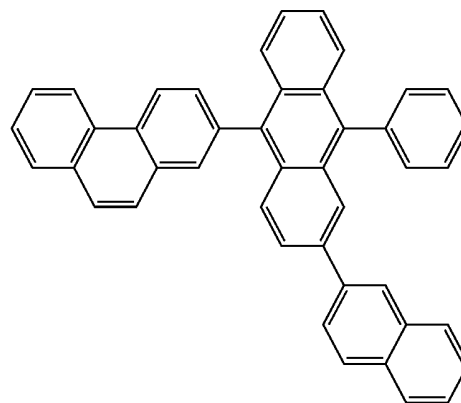


118
-continued

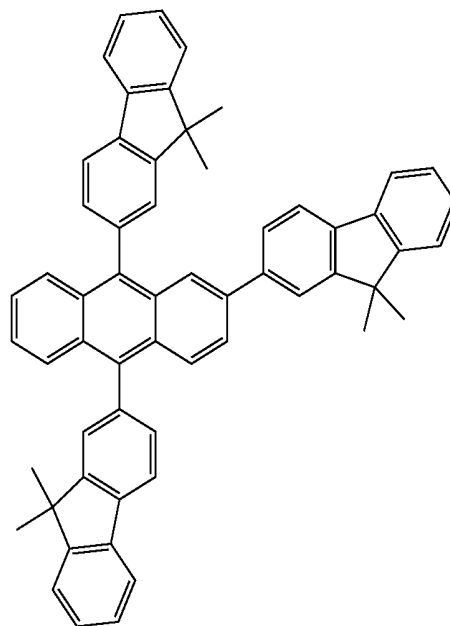
HT23



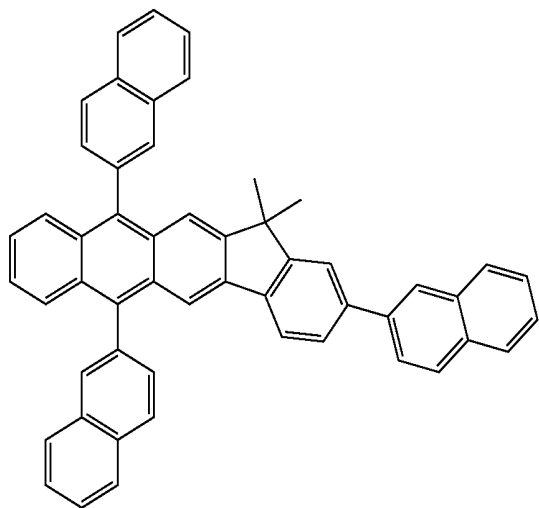
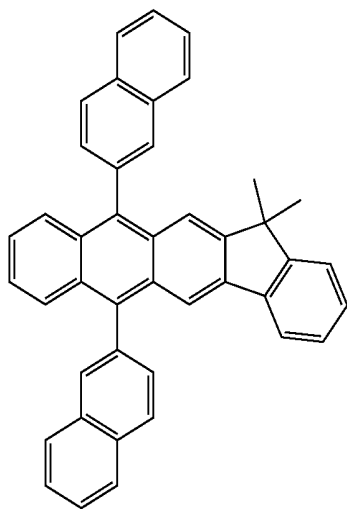
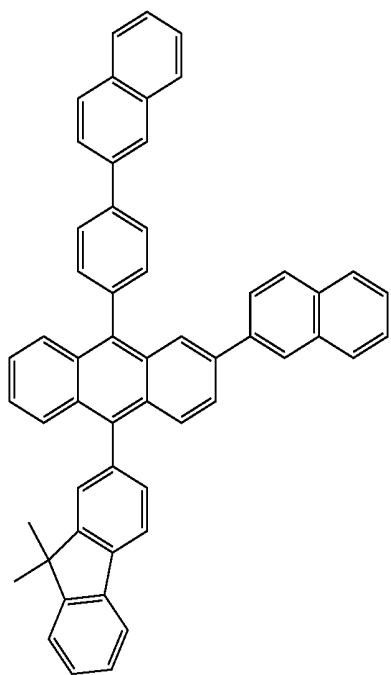
25
H24



45
H25



119
-continued



120
-continued

H26

5

10

15

20

25

H27

30

35

40

H28

45

50

55

60

65

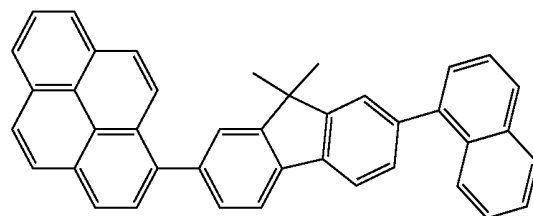
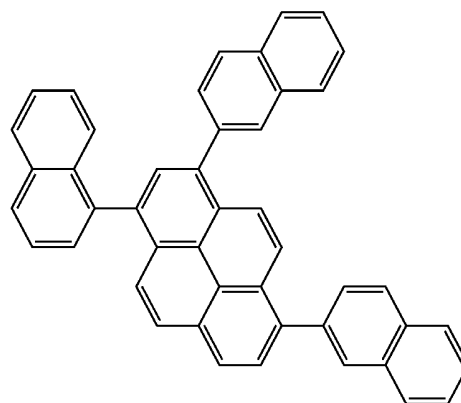
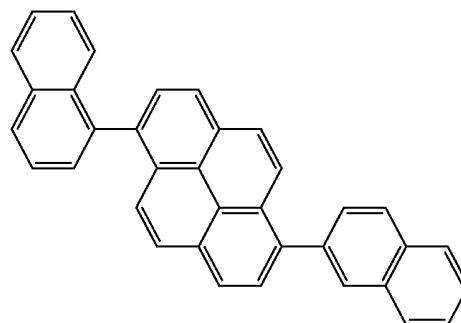
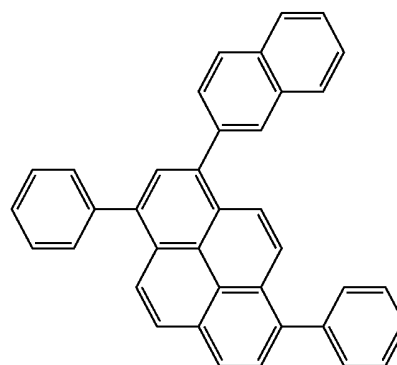
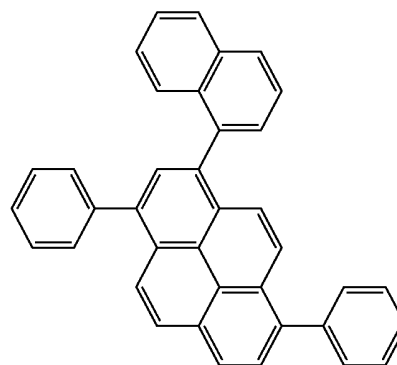
H29

H30

H31

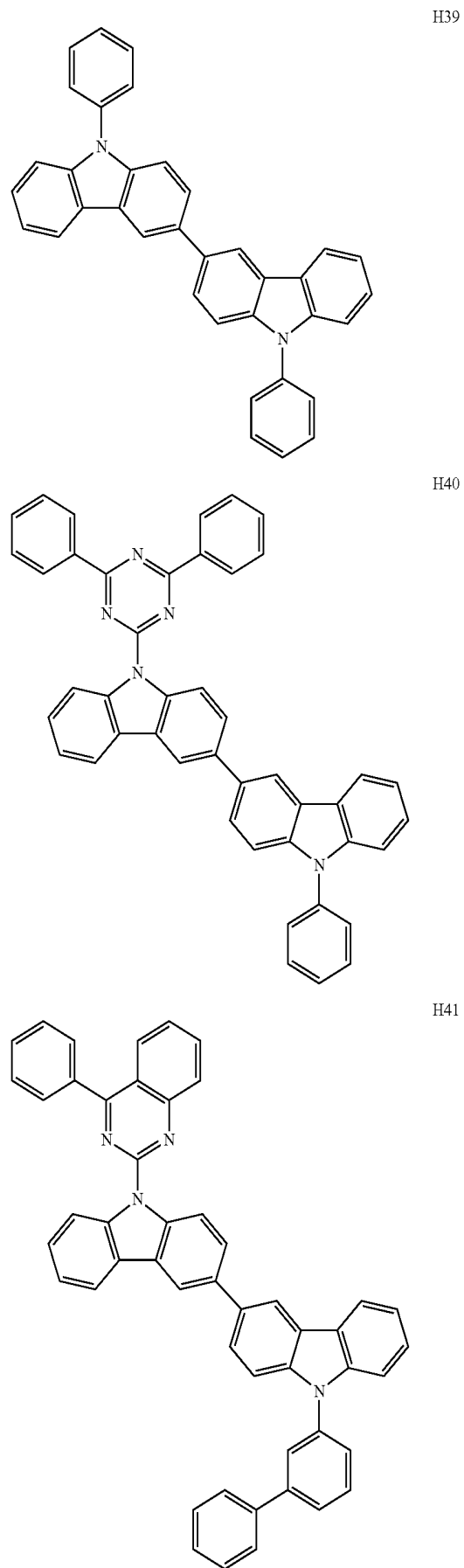
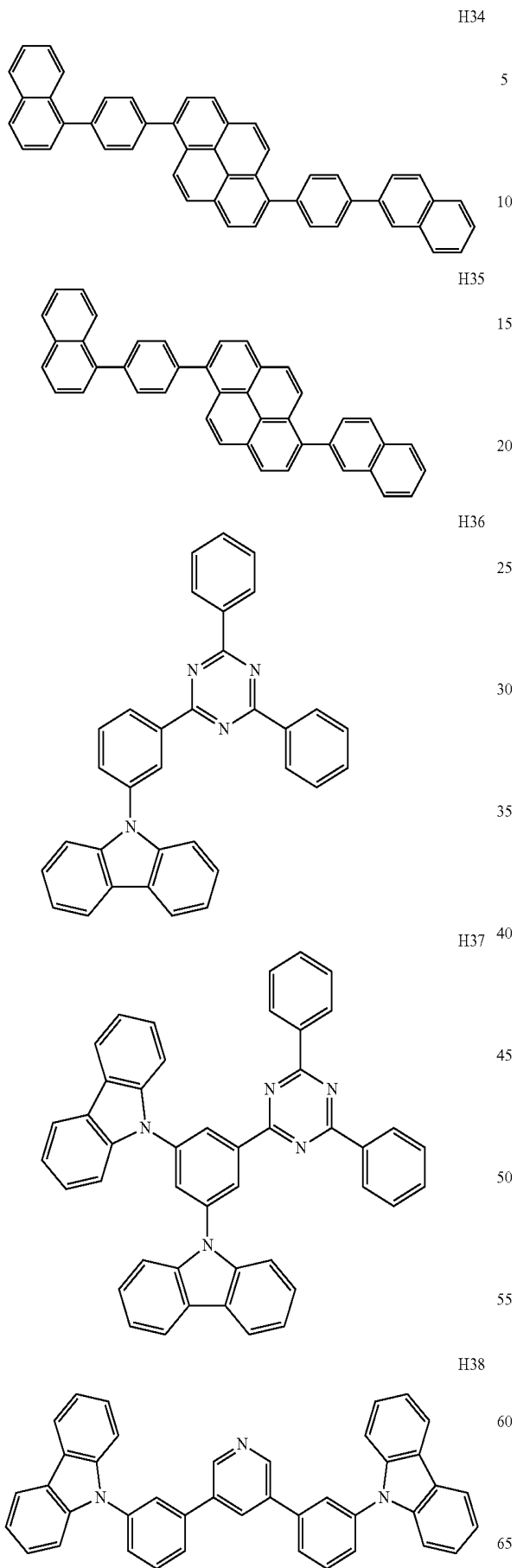
H32

H33

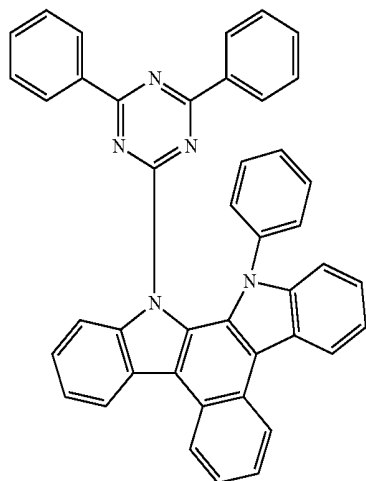
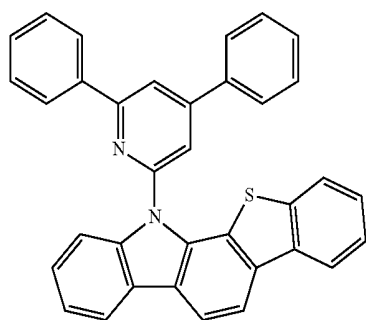
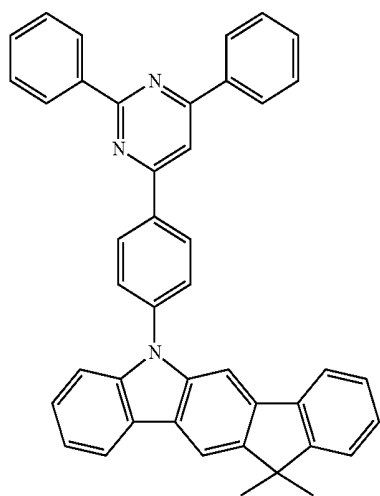
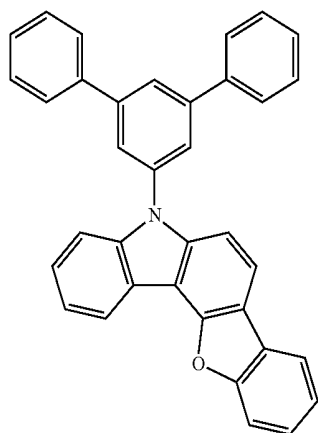


121
-continued

122
-continued



123
-continued



124
-continued

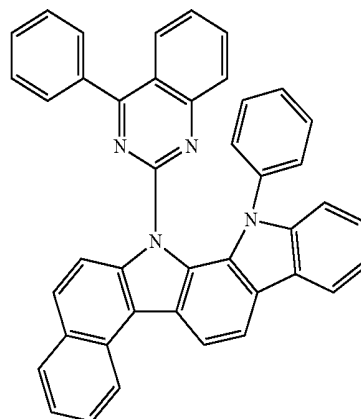
H42

H46

5

10

15

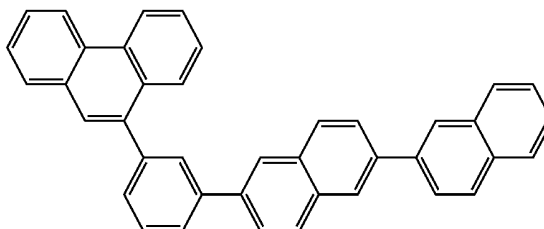


H43

H47

20

25



30

H48

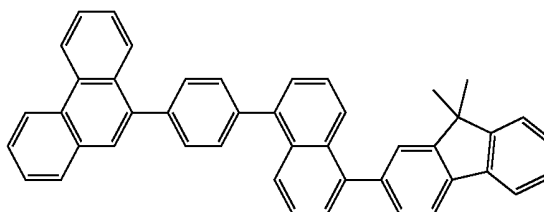
35

H44

H49

40

45

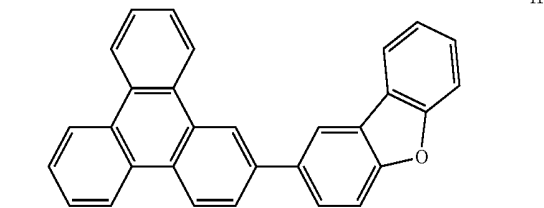


H45

H50

50

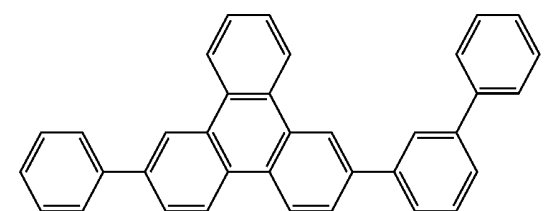
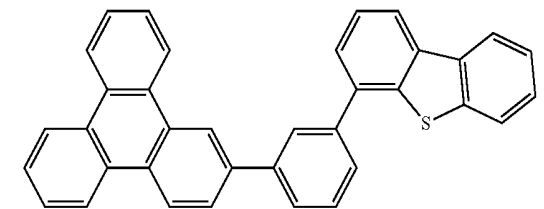
55



60

H51

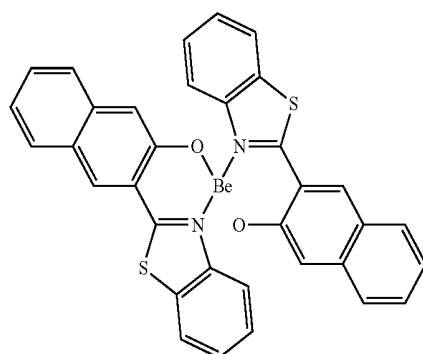
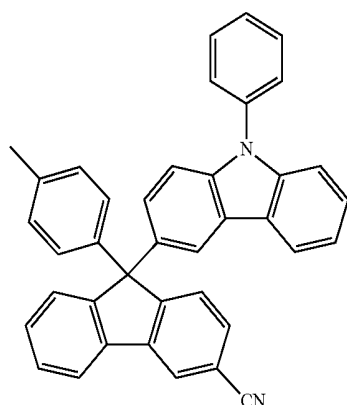
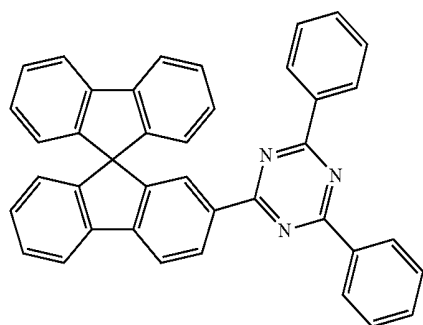
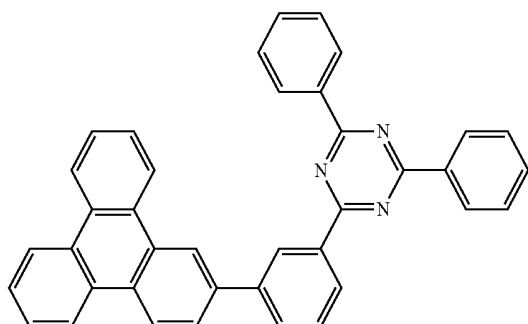
65



125

-continued

H52



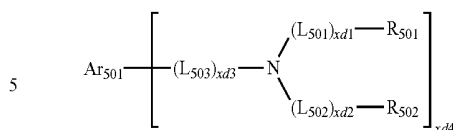
Fluorescent Dopant in Emission Layer

The fluorescent dopant may include an arylamine compound or a styrylamine compound.

The fluorescent dopant may include a compound represented by Formula 501:

126

Formula 501



wherein, in Formula 501,

10 Ar_{501} may be selected from a substituted or unsubstituted $\text{C}_5\text{-C}_{60}$ carbocyclic group and a substituted or unsubstituted $\text{C}_1\text{-C}_{60}$ heterocyclic group,

15 L_{501} to L_{503} may each independently be selected from a substituted or unsubstituted $\text{C}_3\text{-C}_{10}$ cycloalkylene group, a substituted or unsubstituted $\text{C}_1\text{-C}_{10}$ heterocycloalkylene group, a substituted or unsubstituted $\text{C}_3\text{-C}_{10}$ cycloalkenylene group, a substituted or unsubstituted $\text{C}_1\text{-C}_{10}$ heterocycloalkenylene group, a substituted or unsubstituted $\text{C}_6\text{-C}_{60}$ arylene group, a substituted or unsubstituted $\text{C}_1\text{-C}_{60}$ heteroarylene group, a substituted or unsubstituted divalent non-aromatic condensed polycyclic group, and a substituted or unsubstituted divalent non-aromatic condensed heteropolycyclic group,

25 x_{d1} to x_{d3} may each independently be an integer from 0 to 3,

30 R_{501} and R_{502} may each independently be selected from a substituted or unsubstituted $\text{C}_3\text{-C}_{10}$ cycloalkyl group, a substituted or unsubstituted $\text{C}_1\text{-C}_{10}$ heterocycloalkyl group, a substituted or unsubstituted $\text{C}_3\text{-C}_{10}$ cycloalkenyl group, a substituted or unsubstituted $\text{C}_1\text{-C}_{10}$ heterocycloalkenyl group, a substituted or unsubstituted $\text{C}_6\text{-C}_{60}$ aryl group, a substituted or unsubstituted $\text{C}_6\text{-C}_{60}$ aryloxy group, a substituted or unsubstituted $\text{C}_6\text{-C}_{60}$ arylthio group, a substituted or unsubstituted $\text{C}_1\text{-C}_{60}$ heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, and a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, and

x_{d4} may be an integer from 1 to 6.

40 In one embodiment, in Formula 501, Ar_{501} may be selected from

45 a naphthalene group, a heptalene 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, and an indenophenanthrene group; and

50 a naphthalene group, a heptalene 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, and an indenophenanthrene 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 $\text{C}_1\text{-C}_{20}$ alkyl group, a $\text{C}_1\text{-C}_{20}$ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group.

55 In one embodiment, in Formula 501, L_{501} and L_{503} may each independently be selected from

60 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

127

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 carbazolylenylene group, an indolylenylene group, an isoindolylenylene group, a benzofuranylene group, a benzothiophenylene group, a dibenzofuranylene group, a dibenzothiophenylene group, a benzocarbazolylenylene group, a dibenzocarbazolylenylene group, a dibenzosilolylenylene group, and a pyridinylenylene group; and

a phenylene group, a naphthylene group, a fluorenylene group, a spiro-bifluorenylene group, a benzofluorenylene group, a dibenzofluorenylene group, a phenanthrenylene group, an anthracenylenylene 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 carbazolylenylene group, an indolylenylene group, an isoindolylenylene group, a benzofuranylene group, a benzothiophenylene group, a dibenzofuranylene group, a dibenzothiophenylene group, a benzocarbazolylenylene group, a dibenzocarbazolylenylene group, a dibenzosilolylenylene group, and a pyridinylenylene 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 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, and a pyridinyl group.

In one embodiment, in Formula 501, R_{505} and R_{502} may each independently be selected from

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, and a pyridinyl group; and

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, and a pyridinyl 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

128

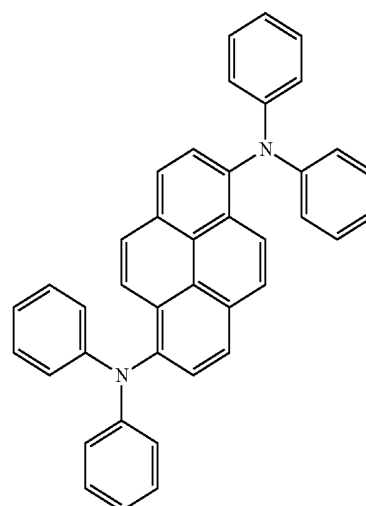
group, a C_1 - C_{20} alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, a fluorenyl group, a spiro-bifluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl 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, and —Si(Q_{31})(Q_{32})(Q_{33}),

wherein Q_{31} to Q_{33} may 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.

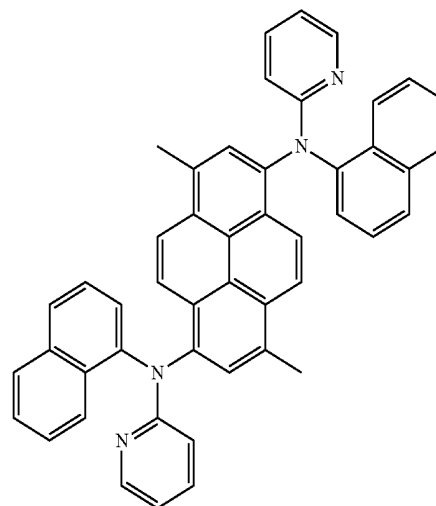
In one or more embodiments, xd4 in Formula 501 may be 2, but embodiments of the present disclosure are not limited thereto.

In some embodiments, the fluorescent dopant may be selected from Compounds FD1 to FD22:

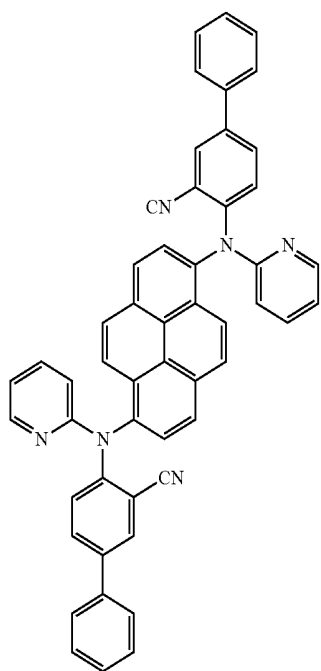
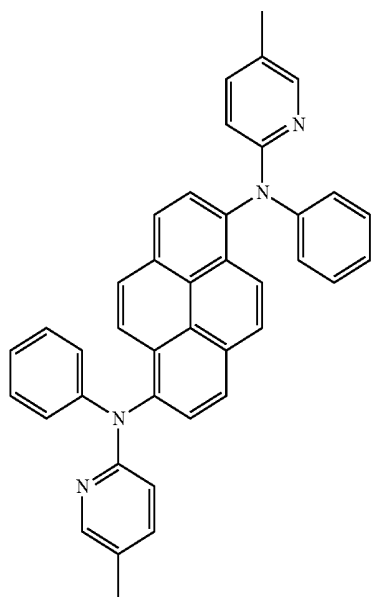
FD1



FD2



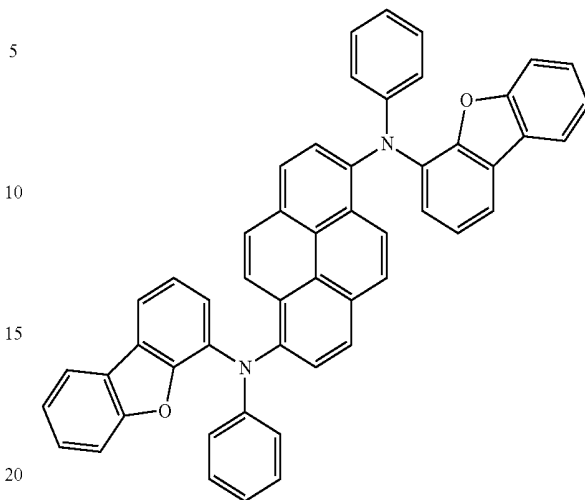
129
-continued



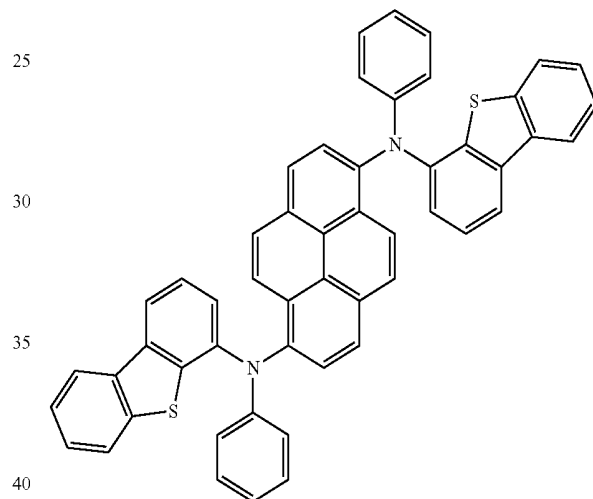
130
-continued

FD3

FD5

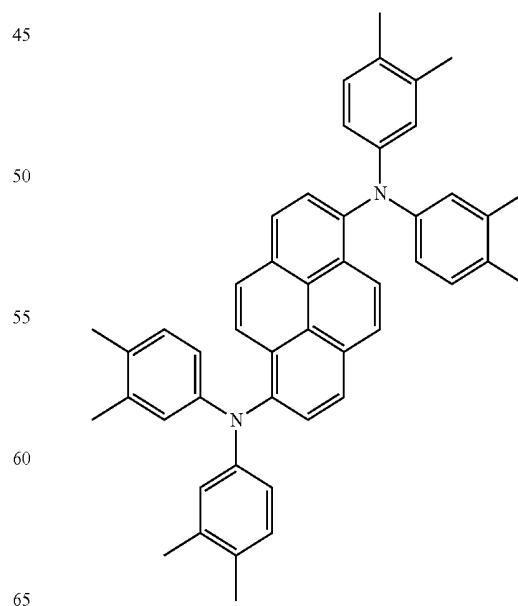


FD6



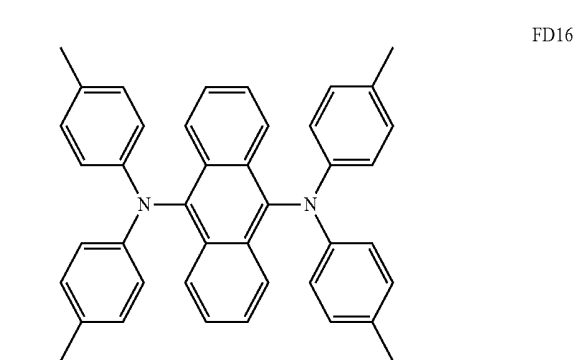
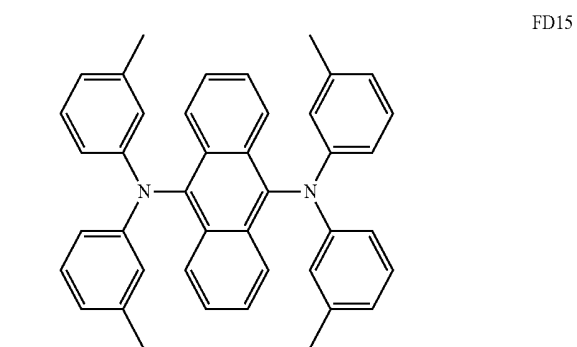
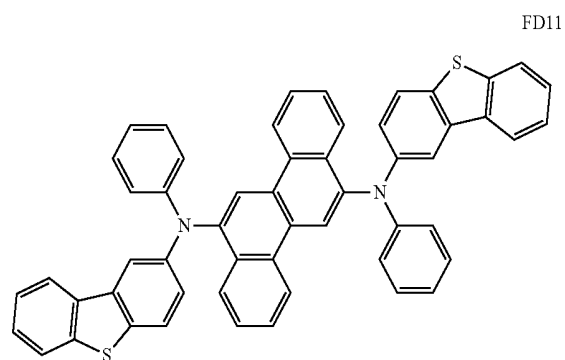
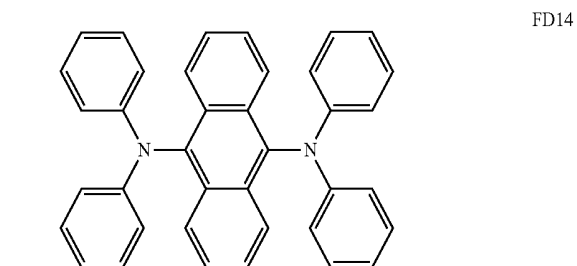
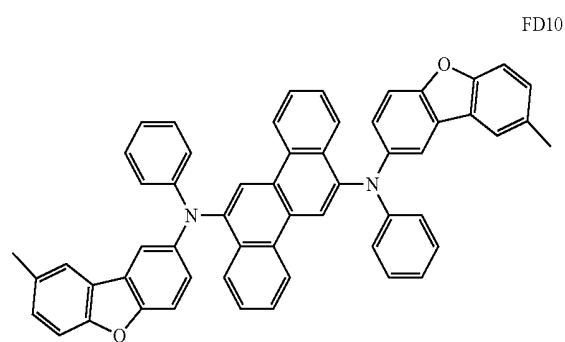
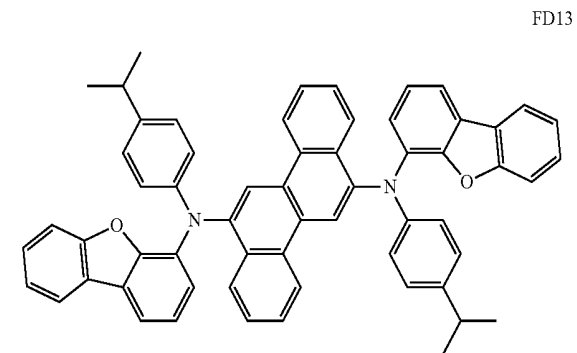
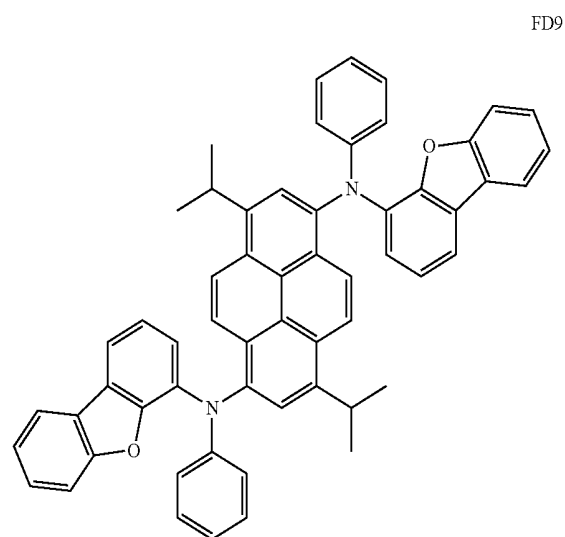
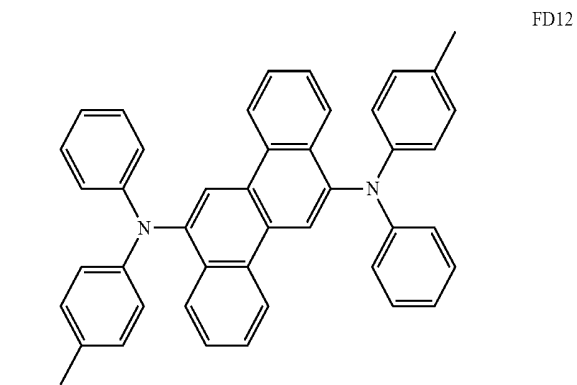
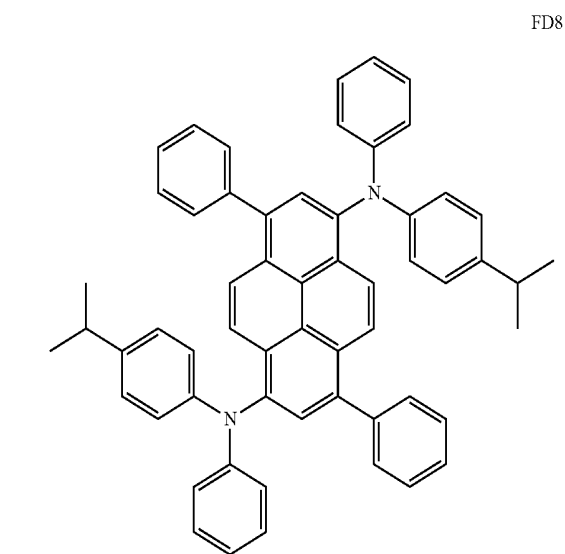
FD4

FD7

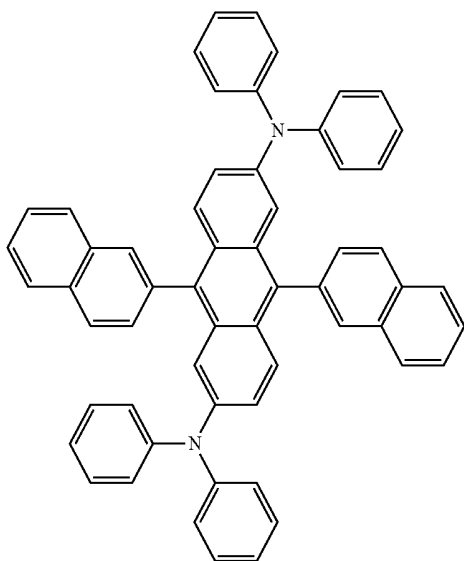
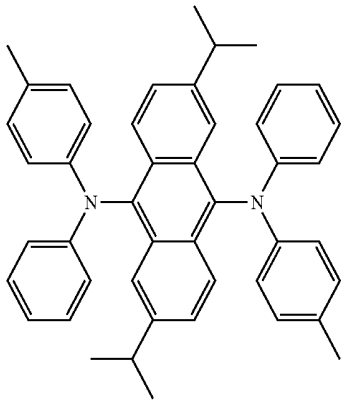
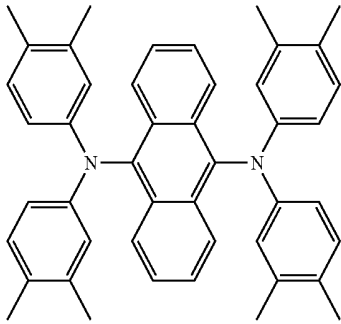


131
-continued

132
-continued



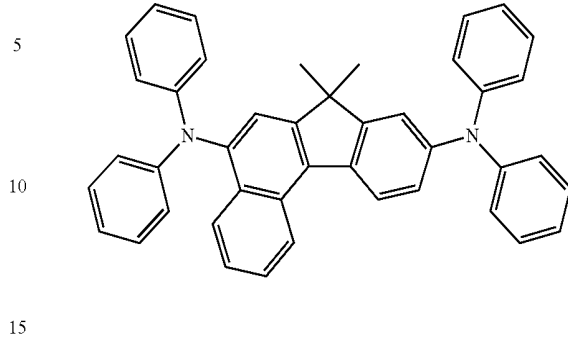
133
-continued



134
-continued

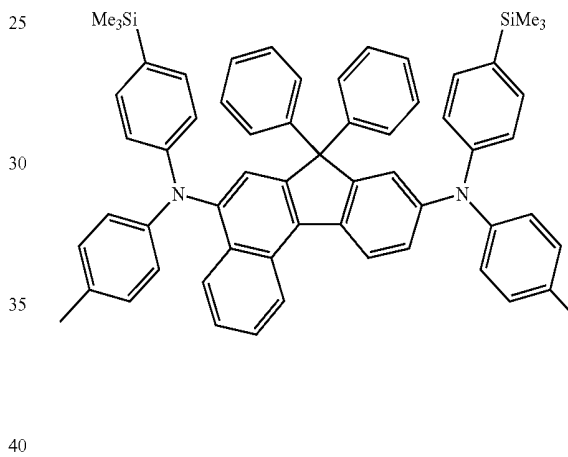
FD17

FD20



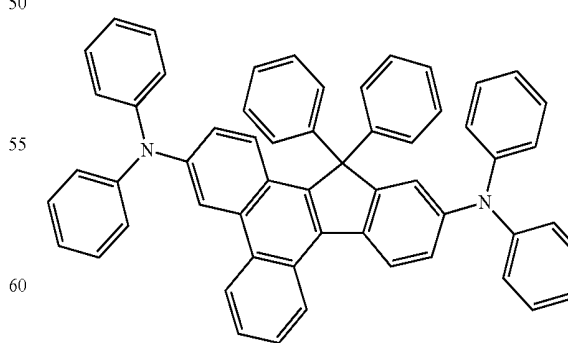
FD18

FD21



FD19

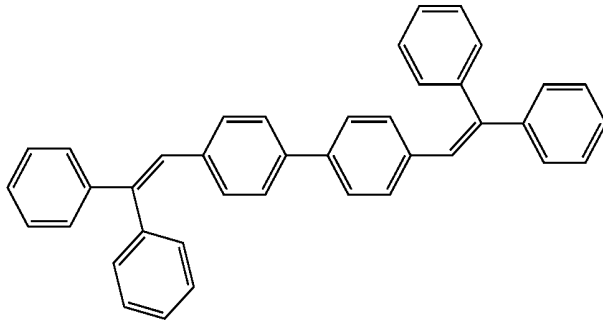
FD22



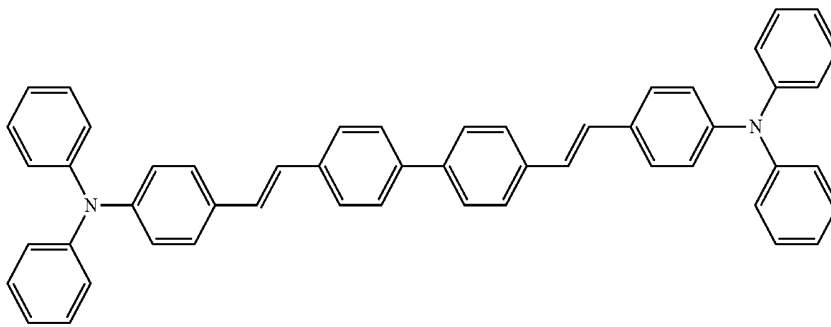
65 In one or more embodiments, the fluorescent dopant may be selected from the following compounds, but embodiments of the present disclosure are not limited thereto:

135

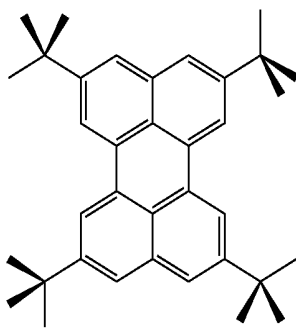
136



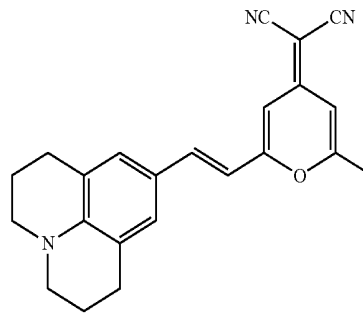
DPVBi



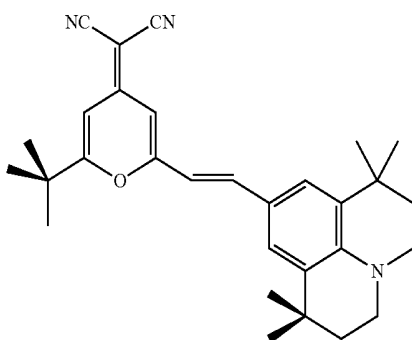
DPAVBi



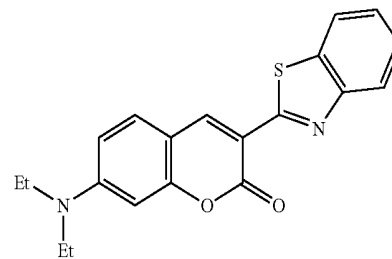
TBPe



DCM

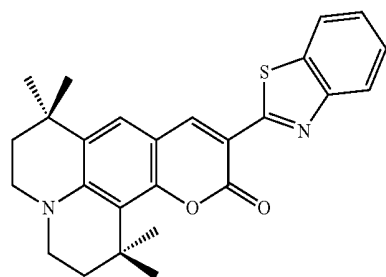


DCJTb



Coumarin 6

-continued



C545T

15

Electron Transport Region in Organic Layer 150

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.

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.

In some embodiments, 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, or an electron control layer/electron transport layer/electron injection layer structure, wherein layers of each structure are sequentially stacked on the emission layer, but embodiments are not limited thereto.

The electron transport region (for example, 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.

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

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, 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.

Examples of the π electron-depleted nitrogen-containing ring may 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, a thiadiazole, an imidazopyridine, an imidazopyrimidine, and an azacarbazole, but embodiments of the present disclosure are not limited thereto.

In some embodiments, the electron transport region may include a compound represented by Formula 601:



wherein, in Formula 601,

Ar_{601} may be selected from a substituted or unsubstituted C_3 - C_{60} carbocyclic group and a substituted or unsubstituted C_1 - C_{60} heterocyclic group,

$xe11$ may be 1, 2, or 3,

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,

$xe1$ may be an integer from 0 to 5,

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 $-\text{P}(=\text{O})(Q_{601})(Q_{602})$,

wherein 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 $xe21$ may be an integer from 1 to 5.

In one embodiment, at least one of the $xe11$ Ar_{601} (s) and the $xe21$ R_{601} (s) may include the foregoing π electron-depleted nitrogen-containing ring.

In one embodiment, in Formula 601, ring Ar_{601} may be selected from

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 iso-benzothiazole group, a benzoxazole group, an isobenzoxazole group, a triazole group, a tetrazole group, an oxadiazole group, a triazine group, a thiaziazole group, an imidazopyridine group, an imidazopyrimidine group, and an azacarbazole group; and

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 iso-benzothiazole group, a benzoxazole group, an isobenzoxazole group, a triazole group, a tetrazole group, an oxadiazole group, a triazine group, a thiaziazole 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),

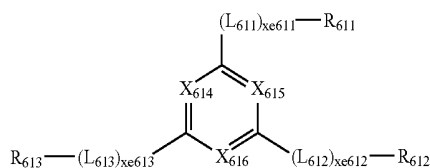
wherein Q₃₁ to Q₃₃ may each independently be selected from a C₁-C₁₀ alkyl group, a C₁-C₁₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group.

When xe11 in Formula 601 is 2 or greater, at least two Ar₆₀₁ (s) may be bound via a single bond.

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

In some embodiments, the compound represented by Formula 601 may be represented by Formula 601-1:

Formula 601-1



wherein, in Formula 601-1,

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

L₆₁₁ to L₆₁₃ may each independently be substantially the same as described above with reference to L₆₀₁,

xe611 to xe613 may each independently be substantially the same as those described herein with reference to xe1,

R₆₁₁ to R₆₁₃ may each independently be substantially the same as described above with reference to R₆₀₁, and

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.

In one embodiment, in Formulae 601 and 601-1, L₆₀₁ and L₆₁₁ to L₆₁₃ may each independently be selected from

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 benzothiothiophenylenylene 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 triazininylenylene group, a quinolininylenylene group, an isoquinolininylenylene group, a benzoquinolininylenylene group, a phthalazininylenylene group, a naphthyridinylenylene group, a quinoxalinylenylene group, a quinazolinylenylene group, a cinnolininylenylene group, a phenanthridinylenylene group, an acridinylenylene group, a phenanthrolinylenylene group, a phenazininylenylene group, a benzimidazolylenylene group, an isobenzothiazolylenylene group, a benzoxazolylenylene group, an isobenzoxazolylenylene group, a triazolylenylene group, a tetrazolylenylene group, an imidazopyridinylenylene group, an imidazopyrimidinylenylene group, and an azacarbazolylenylene group; and

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 benzothiothiophenylenylene 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 triazininylenylene group, a quinolininylenylene group, an isoquinolininylenylene group, a benzoquinolininylenylene group, a phthalazininylenylene group, a naphthyridinylenylene group, a quinoxalinylenylene group, a quinazolinylenylene group, a cinnolininylenylene group, a phenanthridinylenylene group, an acridinylenylene group, a phenanthrolinylenylene group, a phenazininylenylene group, a benzimidazolylenylene group, an isobenzothiazolylenylene group, a benzoxazolylenylene group, an isobenzoxazolylenylene group, a triazolylenylene group, a tetrazolylenylene group, an imidazopyridinylenylene group, an imidazopyrimidinylenylene group, and an

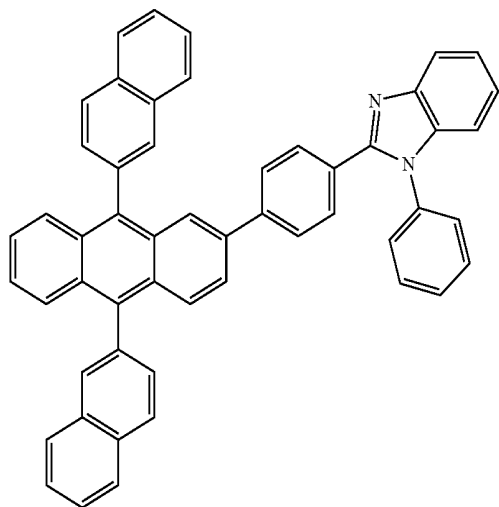
143

144

-continued

ET1

ET4

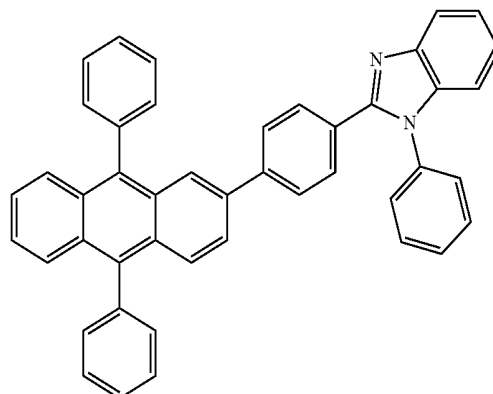


5

10

15

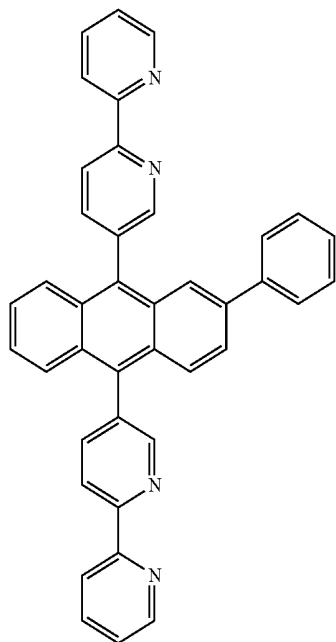
20



ET2

25

ET5

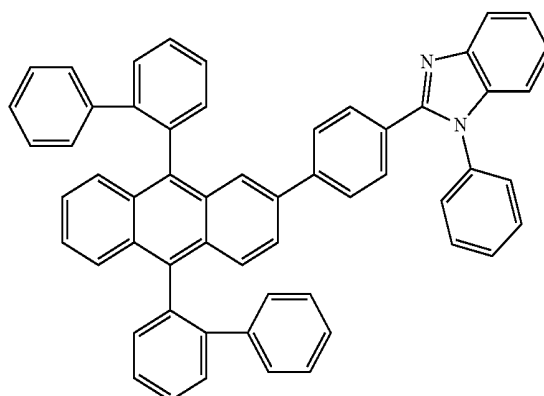


30

35

40

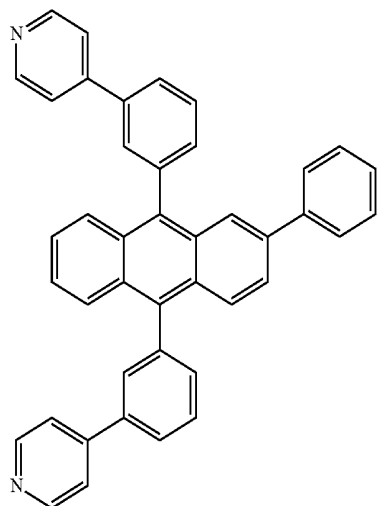
45



ET3

50

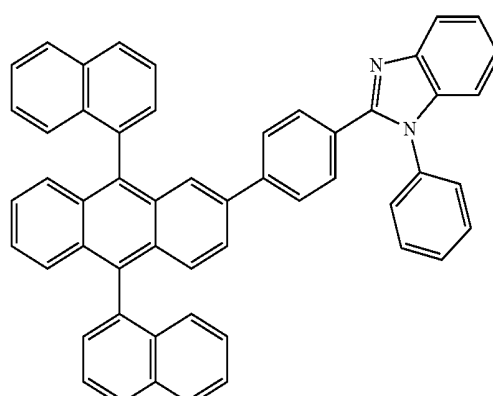
ET6



55

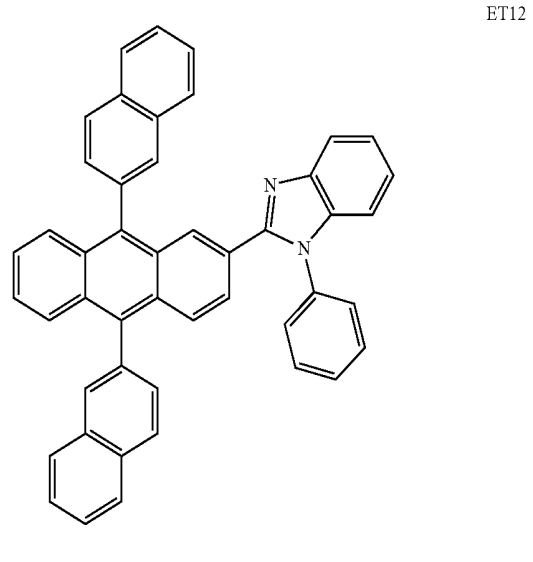
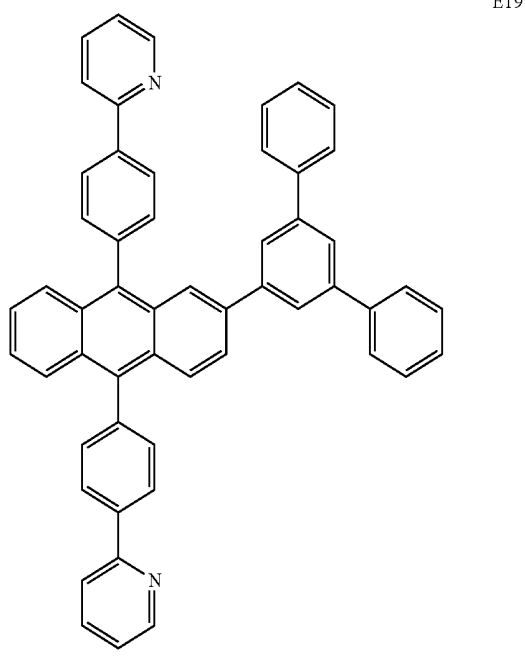
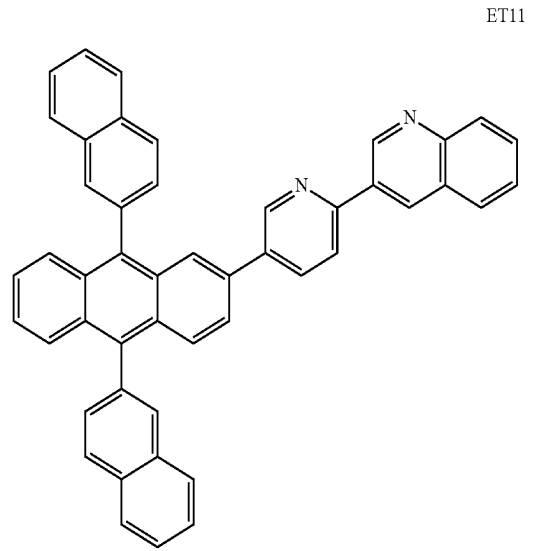
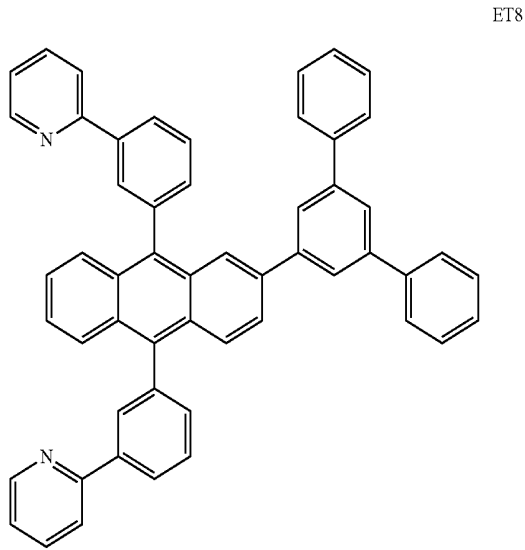
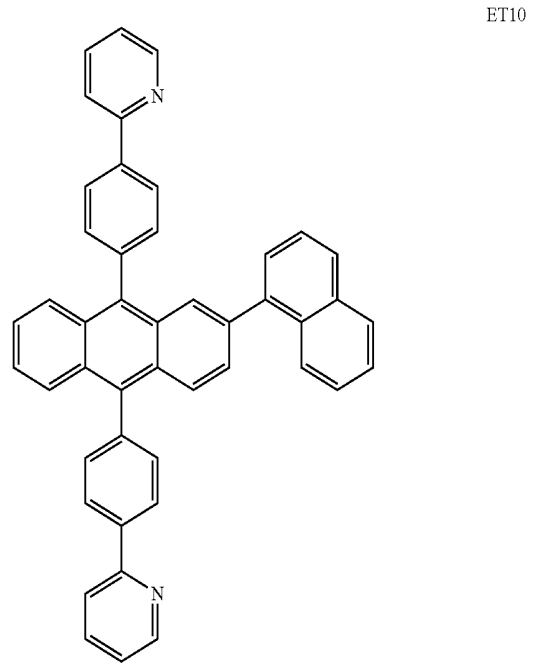
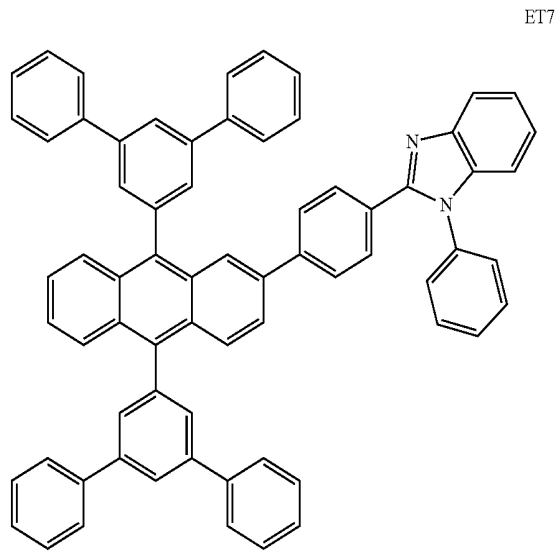
60

65

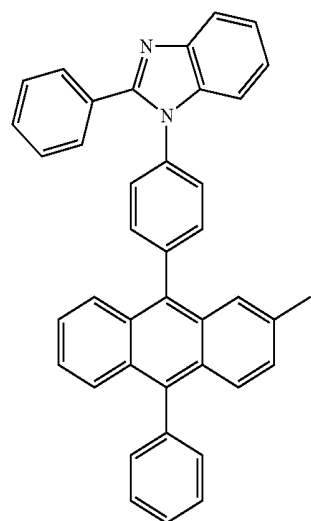
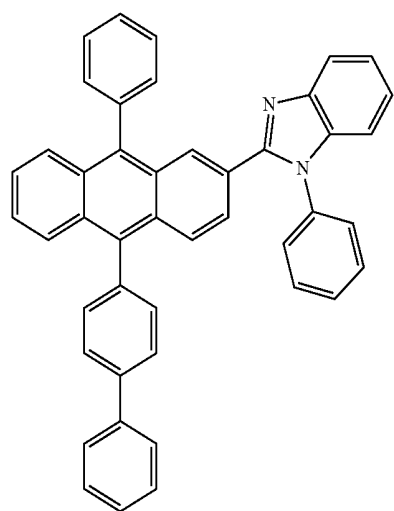
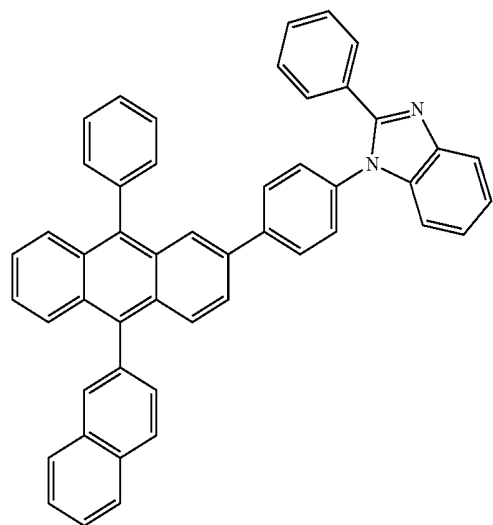


145
-continued

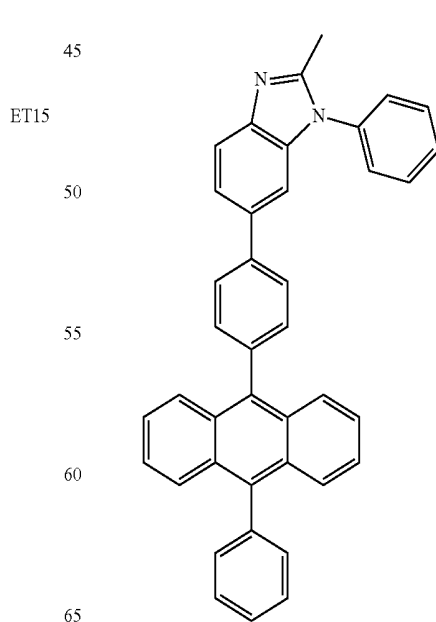
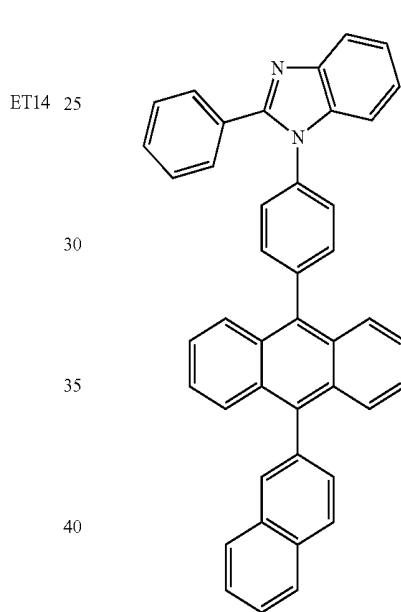
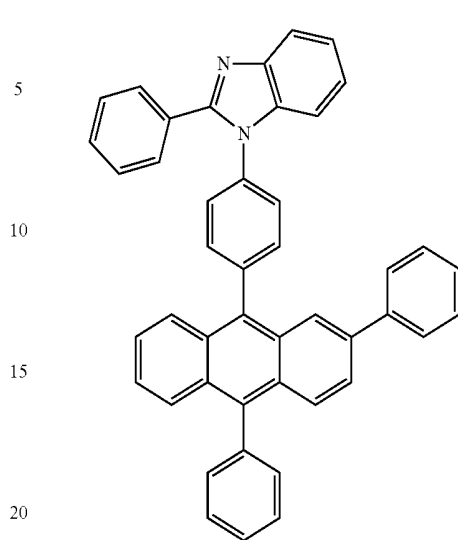
146
-continued



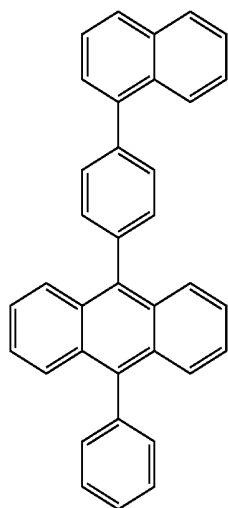
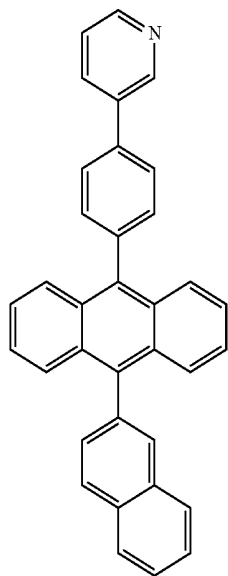
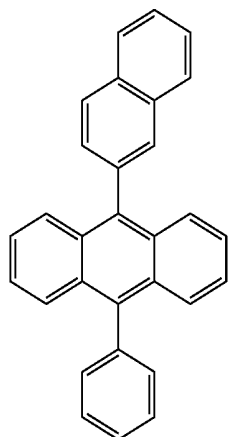
147
-continued



148
-continued



149
-continued



150
-continued

ET19

ET22

5

10

15

20

ET20

25

ET23

30

35

40

45

ET21

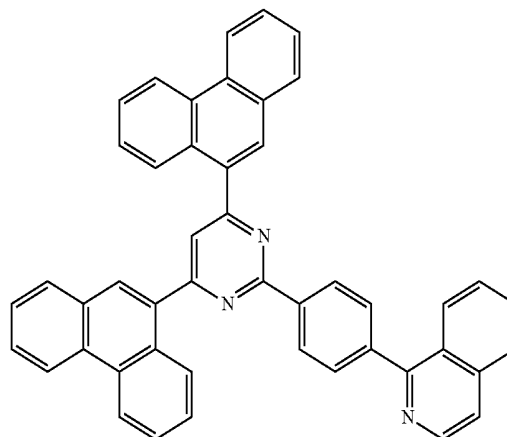
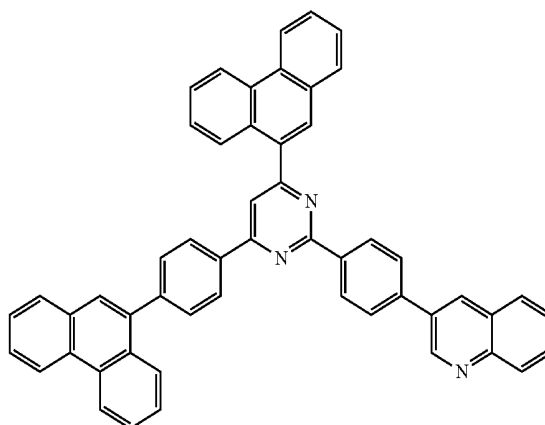
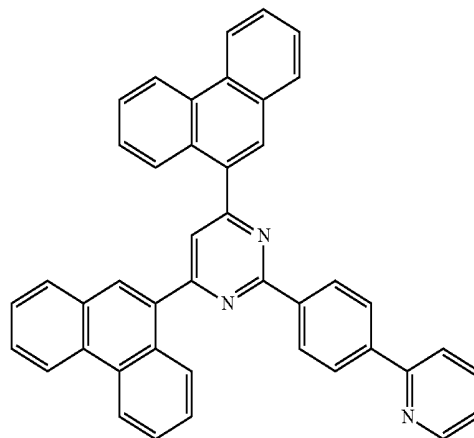
ET24

50

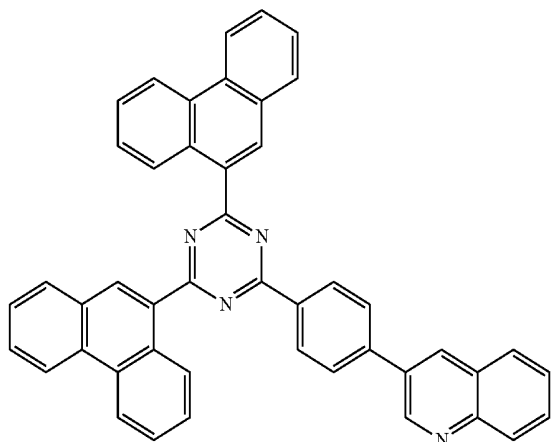
55

60

65



151
-continued



ET25

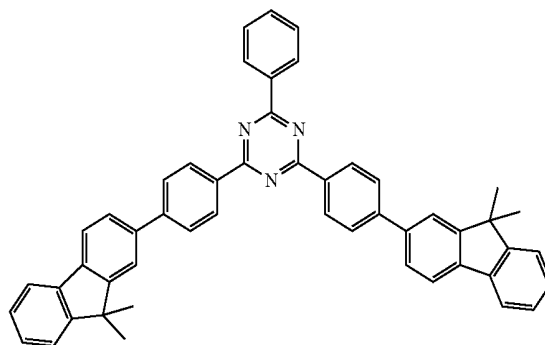
5

10

15

20

152
-continued



ET28

25

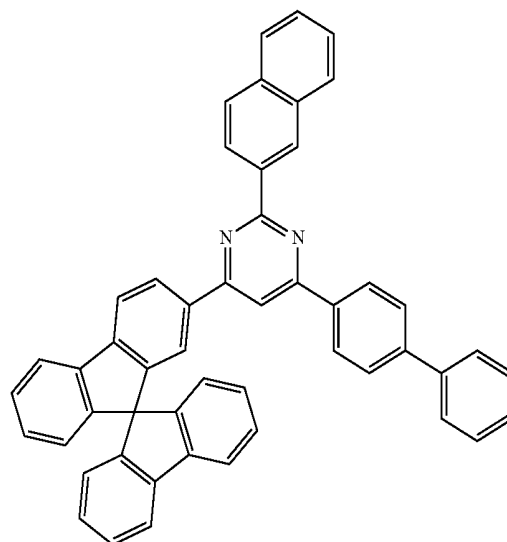
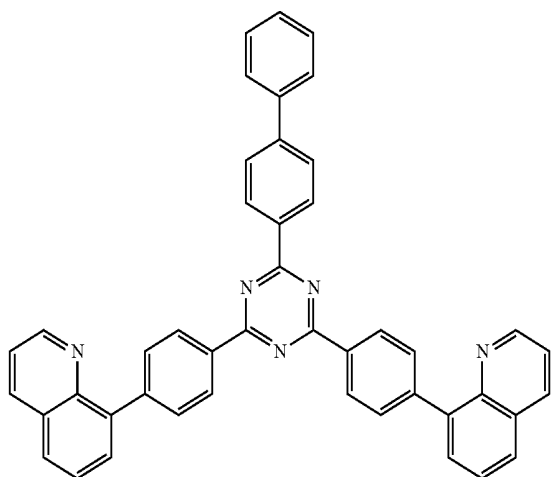
ET26

30

35

40

45



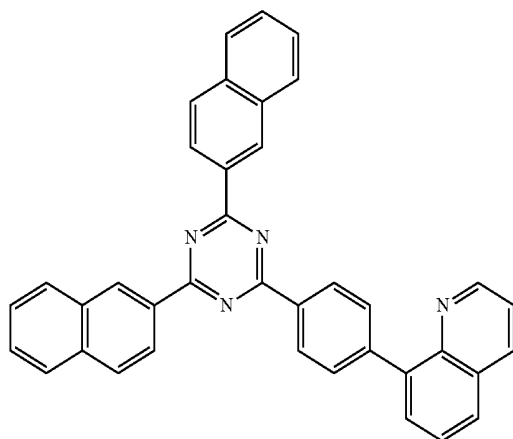
ET27

50

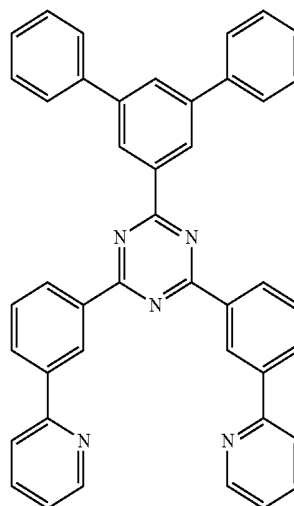
55

60

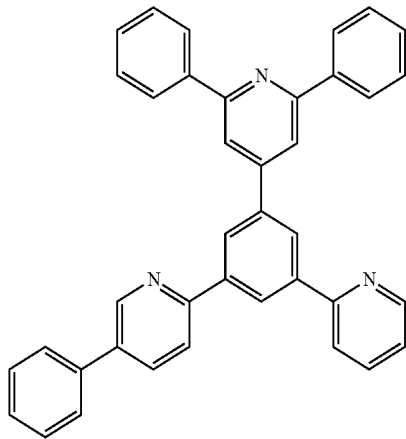
65



ET30



153
-continued



ET31

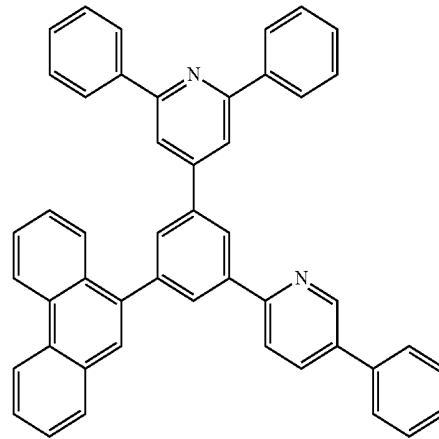
5

10

15

20

154
-continued



ET34

ET32

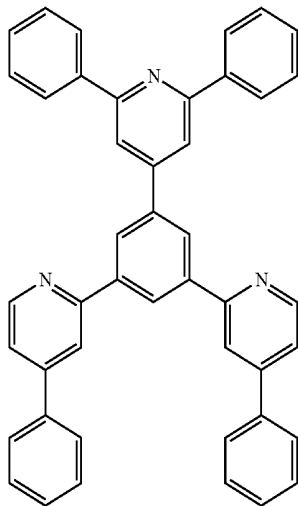
25

30

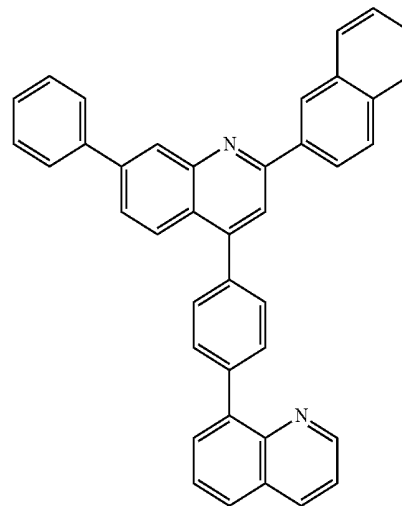
35

40

45



ET35

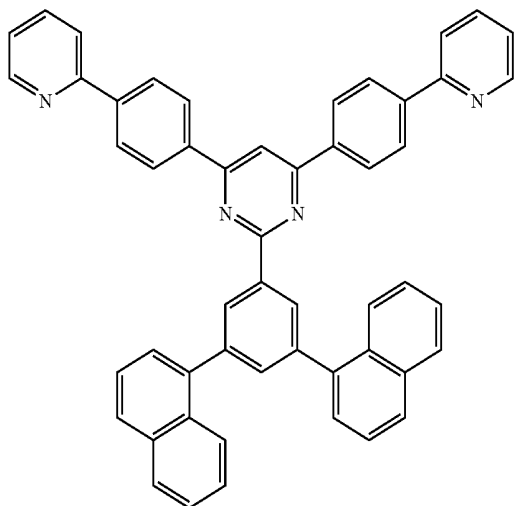


ET33

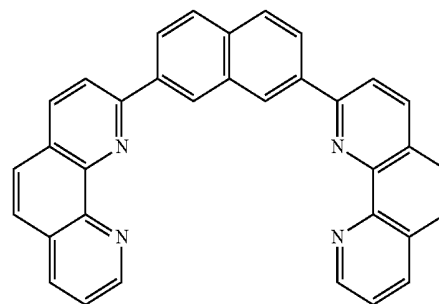
50

55

60

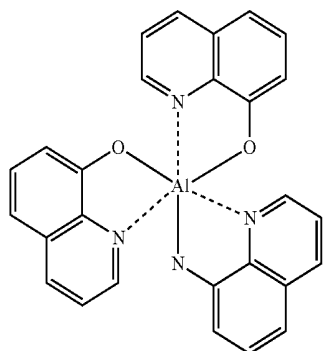
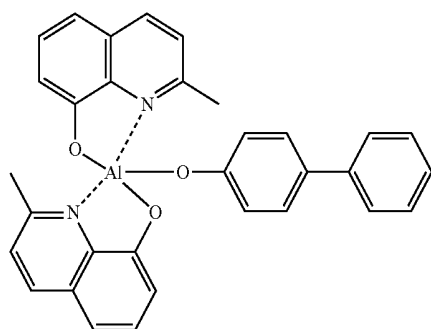


ET36

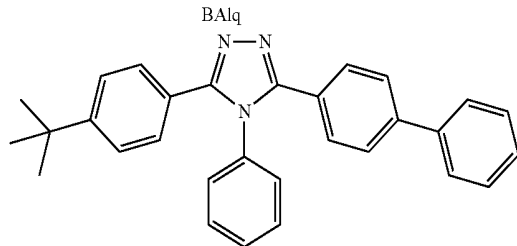


In one or more thickness of the hole blocking layer ors, the electron transport region may include in at one compound selected from 2,9-dimethyl-4,7-diphenyl-1,10-phenanthroline (BCP), 4,7-diphenyl-1,10-phenanthroline (Bphen), Alq3, BAlq, 3-(biphenyl-4-yl)-5-(4-tert-butylphenyl)-4-phenyl-4H-1,2,4-triazole (TAZ), and NTAZ:

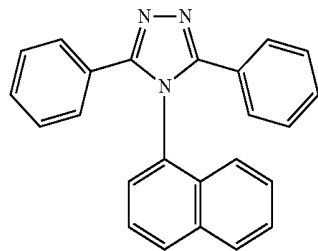
155

Alq₃

BALq



TAZ



NTAZ

The thickness of the hole blocking layer or the electron control layer may be in a range of about 20 Å to about 1,000 Å, and in some embodiments, about 30 Å to about 300 Å. When the thicknesses of the hole blocking layer and the electron control layer are each within any of these ranges, the electron transport region may have excellent electron blocking characteristics or electron control characteristics without a substantial increase in driving voltage.

The thickness of the electron transport layer may be in a range of about 100 Å to about 1,000 Å, and in some embodiments, about 150 Å to about 500 Å. When the thickness of the electron transport layer is within any of these ranges, the electron transport layer may have satisfactory electron transport characteristics without a substantial increase in driving voltage.

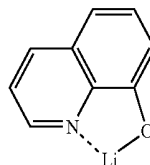
156

The electron transport region (e.g., the electron transport layer in the electron transport region) may further include, in addition to the materials described above, a metal-containing material.

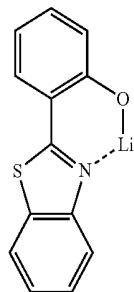
The metal-containing material may include at least one selected from an alkali metal complex and an alkaline earth-metal complex. The alkali metal complex may include a metal ion selected from a Li ion, a Na ion, a K ion, a Rb ion, and a Cs ion. The alkaline earth metal complex may include a metal ion selected from a Be ion, a Mg ion, a Ca ion, a Sr ion, and a Ba ion. Ligands respectively coordinated with the metal ion of the alkali metal complex or the alkaline earth-metal complex may each independently be selected from a hydroxyquinoline, a hydroxyisoquinoline, a hydroxybenzoquinoline, a hydroxyacridine, a hydroxyphenanthridine, a hydroxyphenyl oxazole, a hydroxyphenyl thiazole, a hydroxydiphenyl oxadiazole, a hydroxydiphenyl thiadiazole, a hydroxyphenyl pyridine, a hydroxyphenyl benzimidazole, a hydroxyphenyl benzothiazole, a bipyridine, a phenanthroline, and a cyclopentadiene, but embodiments of the present disclosure are not limited thereto.

In some embodiments, 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



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

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.

The electron injection layer may include at least one selected from 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, and a rare earth metal complex.

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

The alkaline earth-metal may be selected from Ca, Sr, and Ba.

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

The alkali metal compound, the alkaline earth metal compound, and the rare earth metal compound may each independently be selected from oxides and halides (e.g., fluorides, chlorides, bromides, or iodides) of the alkali metal, the alkaline earth metal, and the rare earth metal, respectively.

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 , NaF , CsF , KF , LiI , NaI , CsI , or KI . In one embodiment, the alkali metal compound may be selected from LiF , Li_2O , NaF , LiI , NaI , CsI , and KI , but embodiments of the present disclosure are not limited thereto.

The alkaline earth metal compound may be selected from alkaline earth metal compounds, such as BaO , SrO , CaO , $\text{BaSr}_{1-x}\text{O}$ (wherein $0 < x < 1$), or $\text{BaxCa}_{1-x}\text{O}$ (wherein $0 < x < 1$). In one embodiment, the alkaline earth metal compound may be selected from BaO , SrO , and CaO , but embodiments of the present disclosure are not limited thereto.

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 embodiment, the rare earth metal compound may be selected from YbF_3 , ScF_3 , TbF_3 , YbI_3 , ScI_3 , and TbI_3 , but embodiments are not limited thereto.

The alkali metal complex, the alkaline earth metal complex, and the rare earth metal complex may include an alkali metal ion, an alkaline earth metal ion, and a rare earth metal ion, respectively, as described above, and ligands respectively coordinated with the metal ion of the alkali metal complex, the alkaline earth metal complex, and the rare earth metal complex may each independently be selected from a hydroxyquinoline, a hydroxyisoquinoline, a hydroxybenzoquinoline, a hydroxyacridine, a hydroxyphenanthridine, a hydroxyphenyl oxazole, a hydroxyphenyl thiazole, a hydroxydiphenyl oxadiazole, a hydroxydiphenyl thiadiazole, a hydroxyphenyl pyridine, a hydroxyphenyl benzimidazole, a hydroxyphenyl benzothiazole, a bipyridine, a phenanthroline, and a cyclopentadiene, but embodiments of the present disclosure are not limited thereto.

The electron injection layer may only include the foregoing reducing dopant, or may include the reducing dopant and an organic material. When the electron injection layer includes the reducing dopant and an organic material, the reducing dopant may be homogeneously or un-homogeneously dispersed in a matrix consisting of the organic material.

The thickness of the electron injection layer may be in a range of about 1 Å to about 100 Å, and in some embodiments, about 3 Å to about 90 Å. When the thickness of the electron injection layer is within any of these ranges, the electron injection layer may have satisfactory electron injection characteristics without a substantial increase in driving voltage.

Second Electrode 190

The second electrode 190 may be disposed on the organic layer 150. The second electrode 190 may be a cathode, which is an electron injection electrode, and in this regard, the material for forming the second electrode 190 may be selected from a metal, an alloy, an electrically conductive compound, and a mixture thereof, which may have a relatively low work function.

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

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

Description of FIGS. 2 to 4

An organic light-emitting device 20 illustrated in FIG. 2 includes a first capping layer 210, a first electrode 110, an organic layer 150, and a second electrode 190 which are sequentially stacked in this stated order. An organic light-emitting device 30 illustrated in FIG. 3 includes a first electrode 110, an organic layer 150, a second electrode 190, and a second capping layer 220 which are sequentially stacked in this stated order. An organic light-emitting device 40 illustrated in FIG. 4 includes a first capping layer 210, a first electrode 110, an organic layer 150, a second electrode 190, and a second capping layer 220.

Regarding FIGS. 2 to 4, the first electrode 110, the organic layer 150, and the second electrode 190 may each independently be the same as those described above with reference to FIG. 1.

In the organic layer 150 of each of the organic light-emitting devices 20 and 40, light emitted from the emission layer may pass through the first electrode 110, which is a semi-transmissive electrode or a transmissive electrode, and through the first capping layer 210 toward the outside. In the organic layer 150 of each of the organic light-emitting devices 30 and 40, light emitted from the emission layer may pass through the second electrode 190, which is a semi-transmissive electrode or a transmissive electrode, and through the second capping layer 220 toward the outside.

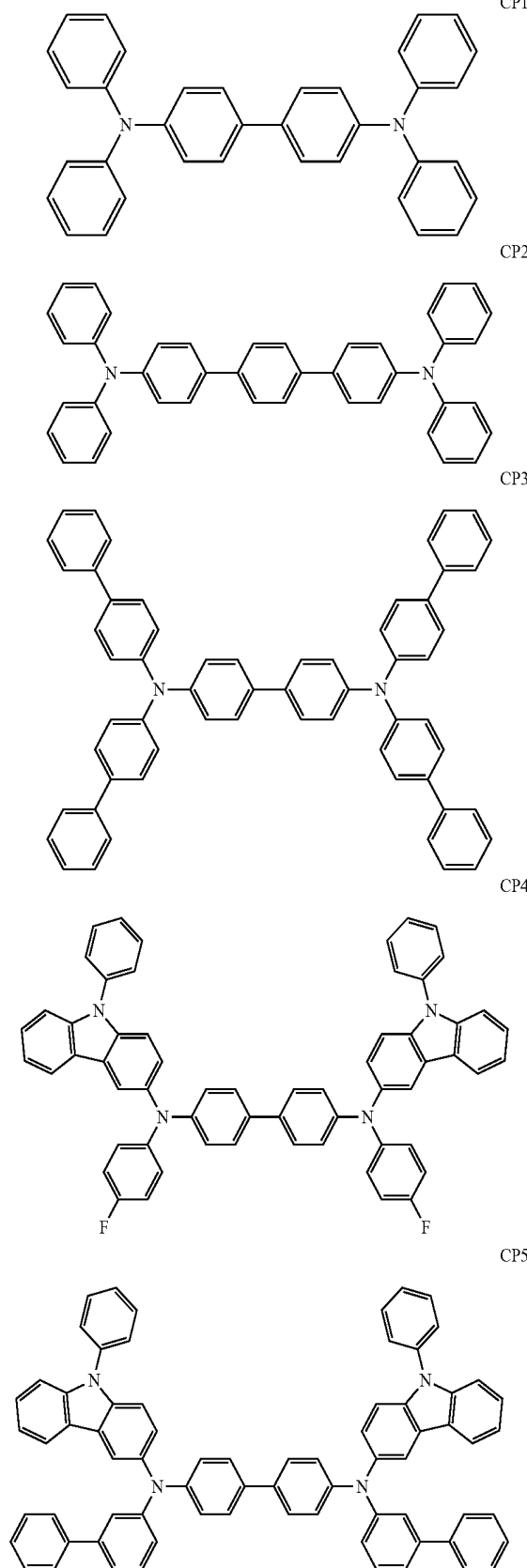
The first capping layer 210 and the second capping layer 220 may increase external luminescent efficiency, based on the principle of constructive interference.

The first capping layer 210 and the second capping layer 220 may each independently be an organic capping layer including an organic material, an inorganic capping layer including an inorganic material, or a composite capping layer including an organic material and an inorganic material.

At least one selected from the first capping layer 210 and the second capping layer 220 may include at least one material selected from carbocyclic compounds, heterocyclic compounds, amine-based compounds, porphyrin derivatives, phthalocyanine derivatives, naphthalocyanine derivatives, alkali metal-based complexes, and alkaline earth metal-based complexes. The carbocyclic compound, the heterocyclic compound, and the amine-based compound may optionally be substituted with a substituent including at least one element selected from oxygen (O), nitrogen (N), sulfur (S), selenium (Se), silicon (Si), fluorine (F), chlorine (Cl), bromine (Br), and iodine (I).

In one embodiment, at least one selected from the first capping layer 210 and the second capping layer 220 may include an amine-based compound.

In one or more embodiments, at least one selected from the first capping layer 210 and the second capping layer 220 may include a compound selected from Compounds CP1 to CP5, but embodiments of the present disclosure are not limited thereto:



Hereinbefore, an organic light-emitting device according to one or more embodiment has been described in connection with FIGS. 1 to 4. However, embodiments are not limited thereto.

The layers constituting the hole transport region, the emission layer, and the layers constituting the electron transport region may be formed in a specific region 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 (LITI).

When the layers constituting the hole transport region, the emission layer, and the layers constituting the electron transport region are each formed by vacuum deposition, for example, vacuum deposition may be performed at a temperature of about 100 to about 500° C., at a vacuum degree of about 10⁻⁸ to about 10⁻³ torr, and at a rate of about 0.01 Angstroms per second (Å/sec) to about 100 Å/sec depending on the compound to be included in each layer and the structure of each layer to be formed.

When the layers constituting the hole transport region, the emission layer, and the layers constituting the electron transport region are each independently formed by spin coating, the spin coating may be performed, for example, at a rate of about 2,000 revolutions per minute (rpm) to about 5,000 rpm and at a heat treatment temperature of about 80° C. to 200° C., depending on the compound to be included in each layer and the structure of each layer to be formed.

25 General Definition of Substituents

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.

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 alkyne group" as used herein refers to a divalent group having the same structure as the C₁-C₆₀ alkyl group.

The term "C₂-C₆₀ alkenyl group" as used herein refers to a hydrocarbon group having at least one carbon-carbon double bond in the middle or at the terminus of the C₂-C₆₀ alkyl group. 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.

The term "C₂-C₆₀ alkynyl group" as used herein refers to a hydrocarbon group having at least one carbon-carbon triple bond in the middle or at the terminus of the C₂-C₆₀ alkyl group. 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.

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). Examples thereof include a methoxy group, an ethoxy group, and an isopropoxy group.

The term "C₃-C₁₀ cycloalkyl group" as used herein refers to a monovalent saturated hydrocarbon monocyclic group having 3 to 10 carbon atoms. 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.

The term "C₁-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. Examples thereof include a 1,2,3,4-oxatriazolidinyl group, a tetrahydrofuran 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.

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 its ring and is not aromatic. 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.

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. Examples of the C₁-C₁₀ heterocycloalkenyl group include a 4,5-dihydro-1,2,3,4-oxatriazolyl group, a 2,3-dihydrofuranlyl 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.

The term “C₆-C₆₀ aryl group” as used herein refers to a monovalent group that has an aromatic system having 6 to 60 carbon atoms. The term “C₆-C₆₀ arylene group” as used herein refers to a divalent group that has an aromatic system having 6 to 60 carbon atoms. 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 independently include two or more rings, the respective rings may be fused.

The term “C₁-C₆₀ heteroaryl group” as used herein refers to a monovalent group having an 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 an 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. 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 independently include two or more rings, the respective rings may be fused.

The term “C₆-C₆₀ aryloxy group” as used herein refers to —OA₁₀₂ (wherein A₁₀₂ is the C₆-C₆₀ aryl group). The term “C₆-C₆₀ arylthio group” as used herein refers to —SA₁₀₃ (wherein A₁₀₃ is the C₆-C₆₀ aryl group).

The term “monovalent non-aromatic condensed polycyclic group” as used herein refers to a monovalent group having two or more rings condensed and only carbon atoms (e.g., 8 to 60 carbon atoms) as ring-forming atoms, wherein the entire molecular structure is non-aromatic. Examples of the monovalent non-aromatic condensed polycyclic group include a fluorenyl group. The term “divalent non-aromatic condensed polycyclic group” as used herein refers to a divalent group having the same structure as the monovalent non-aromatic condensed polycyclic group.

The term “monovalent non-aromatic condensed heteropolycyclic group” as used herein refers to a monovalent group having two or more rings condensed and at least one heteroatom selected from N, O, Si, P, and S, in addition to carbon atoms (e.g., 1 to 60 carbon atoms), as ring-forming atoms, wherein the entire molecular structure is non-aromatic. Examples of the monovalent non-aromatic condensed

heteropolycyclic group may include 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.

The term “C₅-C₆₀ carbocyclic group” as used herein refers to a monocyclic or polycyclic group having 5 to 60 carbon atoms only as ring-forming atoms. The C₅-C₆₀ carbocyclic group may be an aromatic carbocyclic group or a non-aromatic carbocyclic group. The C₅-C₆₀ carbocyclic group may be a ring such as a benzene group, 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 C₅-C₆₀ carbocyclic group, the C₅-C₆₀ carbocyclic group may be a trivalent group or a quadrivalent group.

The term “C₁-C₆₀ heterocyclic group” as used herein refers to a group having substantially the same structure as a C₅-C₆₀ carbocyclic group, with the alteration that at least one heteroatom selected from N, O, Si, P, and S may be used as a ring-forming atom, in addition to carbon atoms (e.g., 1 to 60 carbon atoms).

In the present specification, at least one substituent of the substituted C₅-C₆₀ carbocyclic group, substituted C₁-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, a substituted divalent non-aromatic condensed polycyclic group, a substituted divalent non-aromatic condensed heteropolycyclic group, substituted C₁-C₆₀ alkyl group, substituted C₂-C₆₀ alkenyl group, substituted C₂-C₆₀ alkynyl group, substituted C₁-C₆₀ alkoxy group, substituted C₃-C₁₀ cycloalkyl group, substituted C₁-C₁₀ heterocycloalkyl group, substituted C₃-C₁₀ cycloalkenyl group, substituted C₁-C₁₀ heterocycloalkenyl group, substituted C₆-C₆₀ aryl group, substituted C₆-C₆₀ aryloxy group, substituted C₆-C₆₀ arylthio group, substituted C₁-C₆₀ heteroaryl group, substituted monovalent non-aromatic condensed polycyclic group, and substituted monovalent non-aromatic condensed heteropolycyclic group may be selected from:

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;

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);

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;

163

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),

wherein 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, and a monovalent non-aromatic condensed heteropolycyclic group.

The term “Ph” as used herein represents a phenyl group. The term “Me” as used herein represents a methyl group. The term “Et” as used herein represents an ethyl group. The term “ter-Bu” or “But” as used herein represents a tert-butyl group. The term “OMe” as used herein represents a methoxy group.

The term “biphenyl group” as used therein 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.

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

* and * as used herein, unless defined otherwise, each indicate a binding site to an adjacent atom in the corresponding formula.

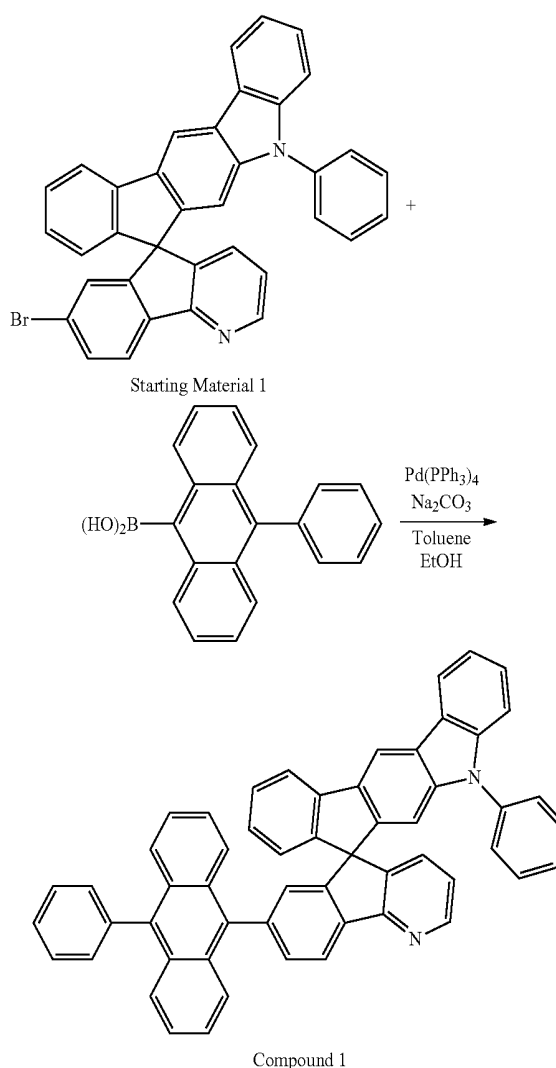
Hereinafter, a compound and an organic light-emitting device, according to one or more embodiments, will be described in detail with reference to Synthesis Examples and Examples. The wording “B was used instead of A” used in describing Synthesis Examples means that an identical molar equivalent of B was used in place of A.

164

EXAMPLES

Synthesis Example

Synthesis Example 1: Synthesis of Compound 1



0.73 g (1 eq, 1.30 mmol) of Starting Material 1 and 0.43 g (1.1 eq, 1.43 mmol) of (10-phenylanthracen-9-yl)boronic acid, and 0.06 g (0.04 eq, 0.052 mmol) of tetrakis(triphenylphosphine)palladium(0) were added to a reaction vessel, and then were dried under vacuum. Next, the reaction vessel was filled with nitrogen gas. 13 ml of toluene was added to the reaction vessel to dissolve the compounds. Then, 6.5 ml of ethanol and 6.5 ml (10 eq, 13.0 mmol) of a 2.0 molar (M) sodium carbonate aqueous solution were added thereto. Then, the mixture was refluxed and stirred at a temperature of 80° C. for 3 hours. Once the reaction was complete, the resultant solution was washed with distilled water, and an organic layer was extracted therefrom using ethyl acetate. The organic layer was dried over magnesium sulfate, filtered by using a Celite filter, and then separated and purified through column chromatography to obtain 0.745 g of Compound 9 (7-(10-phenylanthracen-9-yl)-5'-(1,3,5-triazin-2-yl)-5'H-spiro[indeno[1,2-b]pyridine-5,7'-indeno[2,1-b]carbazole]) (yield: 75%). The obtained compound was identified by mass spectroscopy/fast atom bombardment (MS/FAB).

APCI-MS (m/z): 731[M⁺]

165

Synthesis Example 2: Synthesis of Compound 2

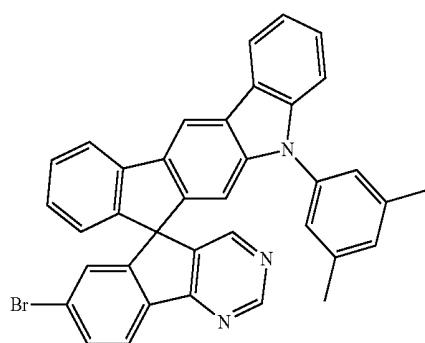
Compound 2 was synthesized in substantially the same manner as in Synthesis Example 1: Synthesis of Compound 1, except that (10-naphthylanthracen-9-yl)boronic acid was used instead of (10-phenylanthracen-9-yl)boronic acid. The obtained compound was identified by MS/FAB.

APCI-MS (m/z): 786[M⁺]

Synthesis Example 3: Synthesis of Compound 17

Compound 17 was synthesized in substantially the same manner as in Synthesis Example 1: Synthesis of Compound 1, except that Starting Material 2 was used instead of Starting Material 1. The obtained compound was identified by ¹H NMR and MS/FAB.

APCI-MS (m/z): 774[M⁺]



Starting Material 2

Example 1

A Corning 15 Ohms per square centimeter (Ω/cm^2 , 1,200 Å) ITO glass substrate was cut to a size of 50 millimeters (mm)×50 mm×0.7 mm, sonicated in isopropyl alcohol and pure water for 5 minutes each, and cleaned by exposure to ultraviolet rays and ozone, so that the resulting glass substrate could be used as a substrate and an anode. Then, the glass substrate was mounted on a vacuum deposition apparatus.

2-TNATA was vacuum-deposited on the ITO anode to form a hole injection layer having a thickness of 600 Å, and 4,4'-bis[N-(1-naphthyl)-N-phenylamino]biphenyl (hereinafter, referred as "NPB") was vacuum-deposited on the hole injection layer to form a hole transport layer having a thickness of 300 Å.

Compound 1 (host) and DPAVBi (dopant) were co-deposited on the hole transport layer at a weight ratio of about 95:5 to form an emission layer having a thickness of about 20 nm.

Compound ET1 was deposited on the emission 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 about 3,000 Å, thereby completing the manufacture of an organic light-emitting device.

166

Examples 2 and 3 and Comparative Examples 1 to 5

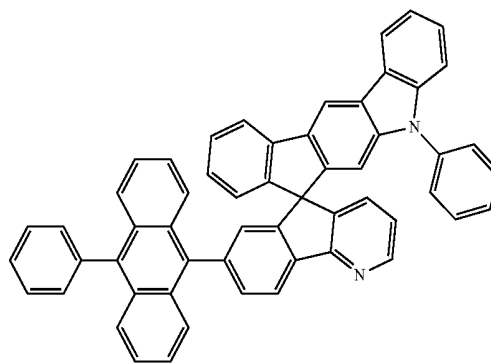
Organic light-emitting devices were manufactured in substantially the same manner as in Example 1, except that the compounds shown in Table 1 were used instead of Compound 1 as a host in forming each emission layer.

Evaluation Example 1

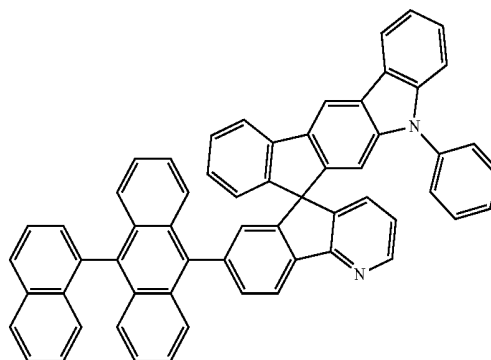
The driving voltage, current density, luminance, and efficiency of the organic light-emitting devices manufactured in Examples 1 to 3 and Comparative Examples 1 to 5 were measured by using a Keithley 236 source-measure unit (SMU) and a PR650 luminance meter. The results thereof are shown in Table 1.

TABLE 1

	Host	Driving voltage (V)	Current density (mA/cm ²)	Luminance (cd/m ²)	Efficiency (cd/A)
25	Example 1 Compound 1	3.65	10	550	5.50
	Example 2 Compound 2	3.7	10	520	5.20
	Example 3 Compound 17	3.2	10	475	4.75
	Comparative Example 1 Compound A	4.6	10	311	3.11
30	Comparative Example 2 Compound B	4.3	10	367	3.67
	Comparative Example 3 Compound C	4.2	10	416	4.16
	Comparative Example 4 Compound D	4.4	10	378	3.78
35	Comparative Example 5 Compound E	4.5	10	252	2.52



1

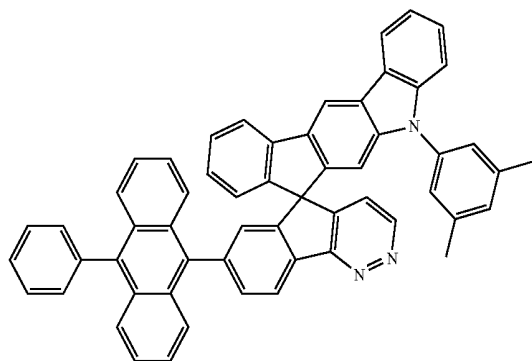


2

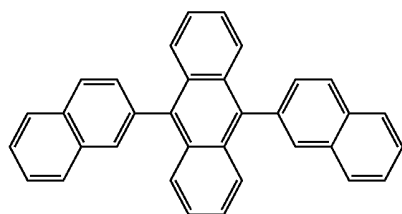
167

TABLE 1-continued

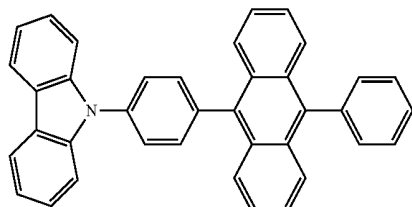
Host	Driving voltage (V)	Current density (mA/cm ²)	Luminance (cd/m ²)	Efficiency (cd/A)
------	---------------------	---------------------------------------	--------------------------------	-------------------



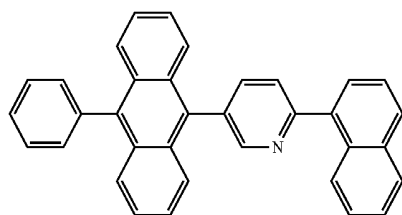
17



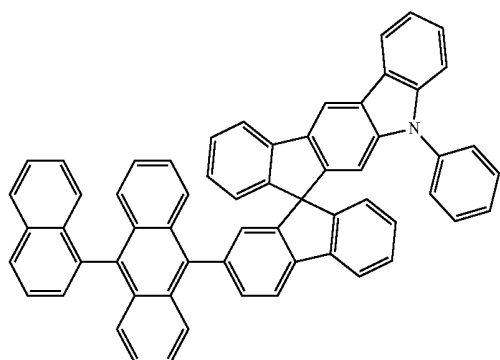
Formula A



Formula B



Formula C

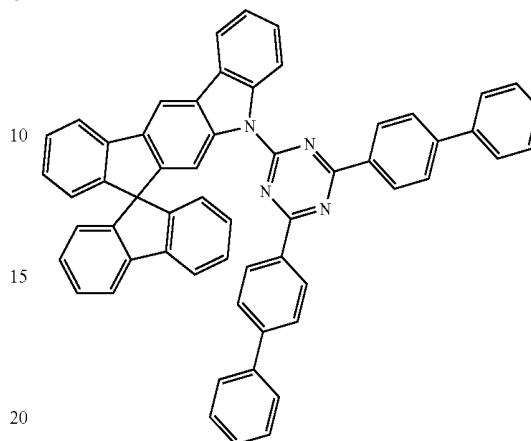


Formula D

168

TABLE 1-continued

Host	Driving voltage (V)	Current density (mA/cm ²)	Luminance (cd/m ²)	Efficiency (cd/A)
------	---------------------	---------------------------------------	--------------------------------	-------------------



Formula E

25 According to the results of Table 1, each of the organic light-emitting devices of Examples 1 to 3 were found to have a low-driving voltage, high luminance, and high efficiency, as compared with the organic light-emitting devices of Comparative Examples 1 to 5.

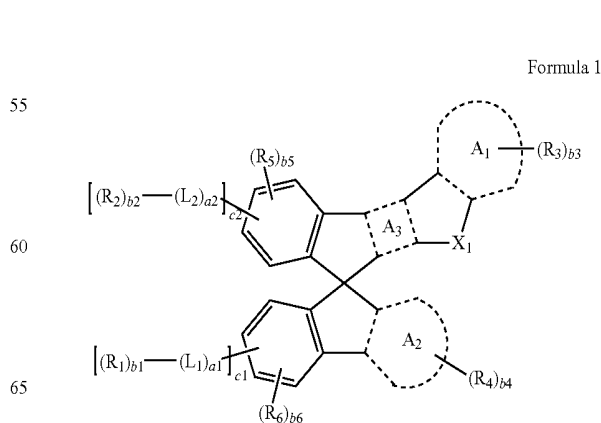
30 As apparent from the foregoing description, an organic light-emitting device including the condensed-cyclic compound may have a low driving voltage, excellent efficiency, and excellent luminance.

35 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.

40 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:

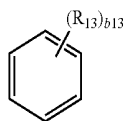
50 1. A condensed-cyclic compound represented by Formula 1:



Formula 1

169

-continued



Formula 2A

5



Formula 2B

10

wherein, in Formulae 1, 2A, and 2B,

ring A_1 is selected from a benzene group, a naphthalene group, a pyridine group, a pyrimidine group, a pyrazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, and a cinnoline group,

ring A_2 is selected from a pyridine group, a pyrimidine group, a pyrazine group, a quinoline group, an isoquinoline group, a quinoxaline group, a quinazoline group, and a cinnoline group,

ring A_3 is selected from groups represented by Formula 2A and groups represented by Formula 2B,

X_1 is selected from N-[(L_{11}) $_{a11}$ -(R_{11}) $_{b11}$], O, and S,

X_2 is selected from N-[(L_{12}) $_{a12}$ -(R_{12}) $_{b12}$], O, and S,

L_1 and L_2 are each independently selected from substituted or unsubstituted condensed polycyclic groups, in which 3 or more carbocyclic groups are condensed with each other,

a_1 and a_2 are each independently an integer from 1 to 5; when a_1 is 2 or greater, at least two L_1 (s) are identical to or different from each other; when a_2 is 2 or greater, at least two L_2 (s) are identical to or different from each other,

L_{11} is selected from a substituted or unsubstituted C_3 - C_{10} cycloalkylene group, a substituted or unsubstituted C_3 - C_{10} cycloalkenylene group, a substituted or unsubstituted C_6 - C_{60} arylene group, and a substituted or unsubstituted divalent non-aromatic condensed polycyclic group,

a_{11} is an integer from 0 to 5,

L_{12} is 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,

a_{12} is an integer from 0 to 5,

R_{11} is selected from 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_6 - C_{60} aryl group, and a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group,

b_{11} is an integer from 1 to 4,

R_1 to R_4 , R_6 , R_{12} , and R_{13} 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

170

substituted or unsubstituted C_1 - C_{60} alkoxy group, a substituted or unsubstituted C_3 - C_{10} cycloalkyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkyl group, a substituted or unsubstituted C_3 - C_{10} cycloalkenyl group, a substituted or unsubstituted C_1 - C_{10} heterocycloalkenyl group, a substituted or unsubstituted C_6 - C_{60} aryl group, a substituted or unsubstituted C_6 - C_{60} aryloxy group, a substituted or unsubstituted C_6 - C_{60} arylthio group, a substituted or unsubstituted C_1 - C_{60} heteroaryl group, a substituted or unsubstituted monovalent non-aromatic condensed polycyclic group, a substituted or unsubstituted monovalent non-aromatic condensed heteropolycyclic group, Si(Q_1)(Q_2)(Q_3), —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_5 is 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 C_1 - C_{20} alkyl group, a C_1 - C_{20} alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group,

b_1 , b_2 , b_5 , b_6 , and b_{12} are each independently an integer from 0 to 4,

b_3 and b_4 are each independently an integer from 0 to 6,

b_{13} is selected from 0, 1, and 2,

c_1 is an integer from 1 to 4 and c_2 is 0,

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

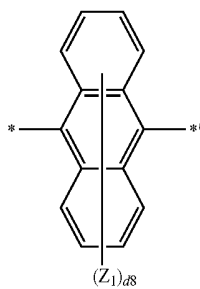
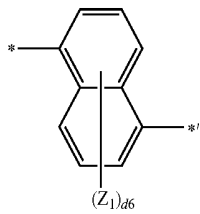
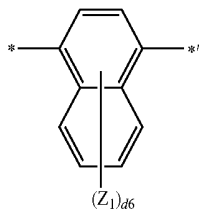
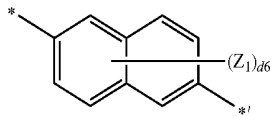
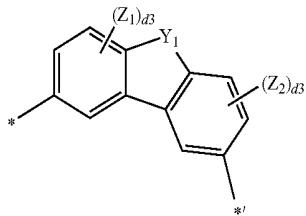
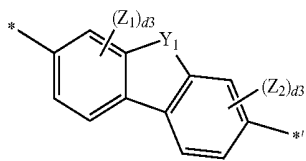
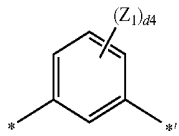
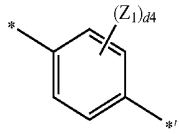
deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group;

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} hetero-

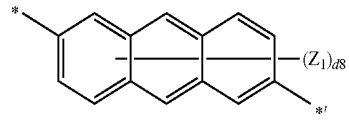
- cycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, a C₁-C₆₀ heteroaryl group, a monovalent non-aromatic condensed polycyclic group, a monovalent non-aromatic condensed heteropolycyclic group, a biphenyl group, and a terphenyl group;
- a C₃-C₁₀ cycloalkyl group, 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),
- wherein 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 C₆-C₆₀ aryl group substituted with a C₁-C₆₀ alkyl group, a C₆-C₆₀ aryl group substituted with a C₆-C₆₀ aryl group, a terphenyl group, a C₁-C₆₀ heteroaryl group, a C₁-C₆₀ heteroaryl group substituted with a C₁-C₆₀ alkyl group, a C₁-C₆₀ heteroaryl group substituted with a C₆-C₆₀ aryl group, a monovalent non-aromatic condensed polycyclic group, and a monovalent non-aromatic condensed heteropolycyclic group.
2. The condensed-cyclic compound of claim 1, wherein ring A₁ is selected from a benzene group and a naphthalene group, and ring A₂ is selected from a pyridine group, a quinoline group, and an isoquinoline group.
3. The condensed-cyclic compound of claim 1, wherein X₁ is N-[(L₁₁)_{a11}-(R₁₁)_{b11}].
4. The condensed-cyclic compound of claim 1, wherein L₁ to L₂ are each independently selected from an indacenylene group, an acenaphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzo-fluorenylene group, a dibenzofluorenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylene group, a fluoranthrenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a naphthacenylene group, a picenylene group, a perylenylene group, a pentaphenylene group, a hexacenylene group, a pentacenylene group, a rubicenylene group, a coronenylene group, and an ovalenylene group; and

- an indacenylene group, an acenaphthylene group, a fluorenylene group, a spiro-fluorenylene group, a benzo-fluorenylene group, a dibenzofluorenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylene group, a fluoranthrenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylene group, a naphthacenylene group, a picenylene group, a perylenylene group, a pentaphenylene group, a hexacenylene group, a pentacenylene group, a rubicenylene group, a coronenylene group, and an ovalenylene 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 cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, 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-fluorenyl group, a benzo-fluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthrenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a 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, a thiadiazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and —Si(Q₃₁)(Q₃₂)(Q₃₃),
- wherein Q₃₁ to Q₃₃ are each independently 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.
5. The condensed-cyclic compound of claim 1, wherein L₁ and L₂ are each independently selected from groups represented by Formulae 3-8, 3-9, 3-25, and 3-35 to 3-41,
- L₁₁ is selected from groups represented by Formulae 3-1 to 3-9, 3-25, and 3-33 to 3-41, wherein Y₁ in Formulae 3-3 and 3-4 is C(Z₃)(Z₄), and
- L₁₂ is selected from groups represented by Formulae 3-1 to 3-41:



Formula 3-1

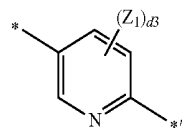
5



Formula 3-9

Formula 3-2

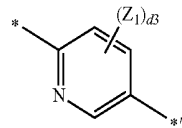
10



Formula 3-10

Formula 3-3

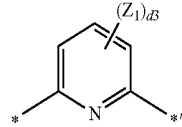
15



Formula 3-11

Formula 3-4

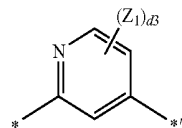
20



Formula 3-12

Formula 3-5

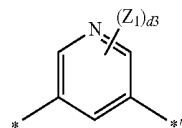
25



Formula 3-13

Formula 3-6

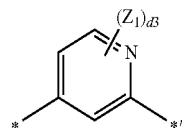
30



Formula 3-14

Formula 3-7

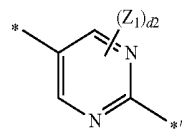
35



Formula 3-15

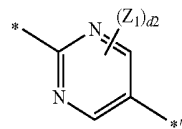
Formula 3-8

40



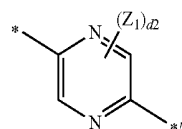
Formula 3-16

45



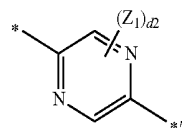
Formula 3-17

50



Formula 3-18

55

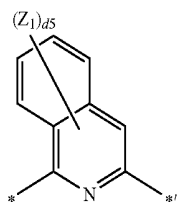
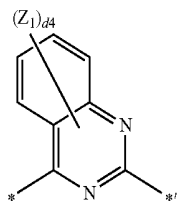
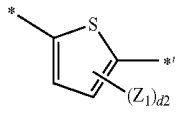
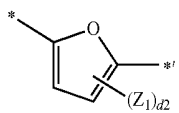
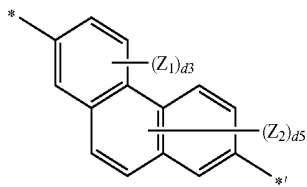
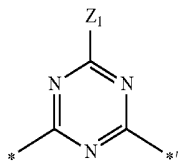
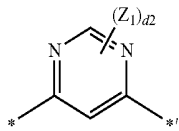
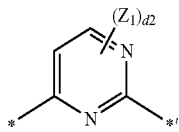
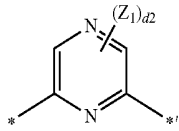
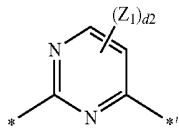


Formula 3-19

60

65

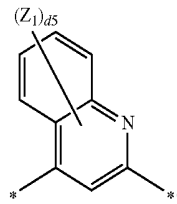
175
-continued



176
-continued

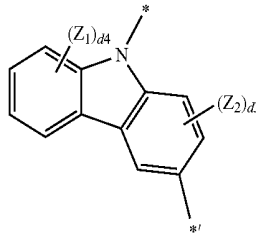
Formula 3-20

5



Formula 3-21

10

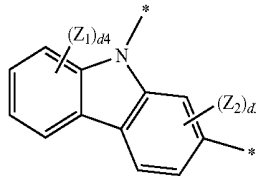


Formula 3-22

15

Formula 3-23

20

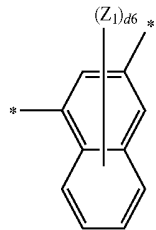


Formula 3-24

25

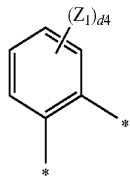
Formula 3-25

35



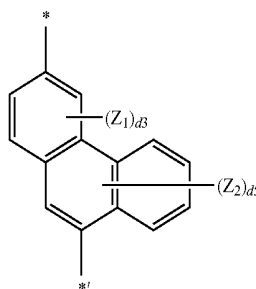
Formula 3-26

40



Formula 3-27

45



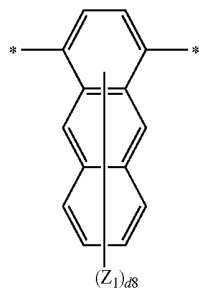
Formula 3-28

50

55

Formula 3-29

60



65

Formula 3-30

Formula 3-31

Formula 3-32

Formula 3-33

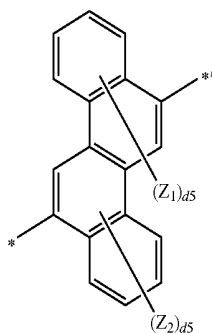
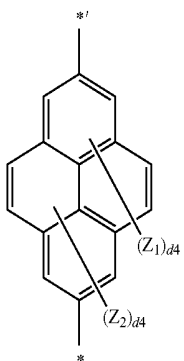
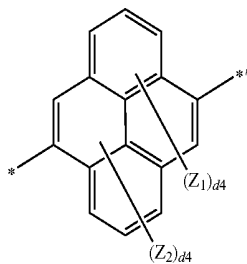
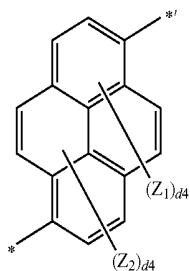
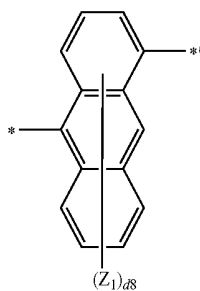
Formula 3-34

Formula 3-35

Formula 3-36

177

-continued



wherein, in Formulae 3-1 to 3-41,
 Y₁ is selected from O, S, C(Z₃)(Z₄), N(Z₅), and Si(Z₆)(Z₇),
 Z₁ to Z₇ are each independently selected from hydrogen, 65
 deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a
 cyano group, a nitro group, an amidino group, a

178

Formula 3-37

5

10

Formula 3-38

15

20

Formula 3-39

25

30

Formula 3-40

35

40

45

Formula 3-41

50

55

60

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 pyrenyl group, a chrysenyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a dibenzofuran group, a dibenzothiophenyl group, a triazinyl group, a benzimidazolyl group, a phenanthrolinyl group, and —Si(Q₃₁)(Q₃₂)(Q₃₃),

wherein Q₃₁ to Q₃₃ are each independently 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,

d2 is an integer selected from 1 and 2,

d3 is an integer from 1 to 3,

d4 is an integer from 1 to 4,

d5 is an integer from 1 to 5,

d6 is an integer from 1 to 6,

d8 is an integer from 1 to 8, and

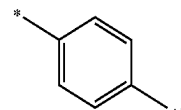
* and *¹ each indicate a binding site to an adjacent atom.

6. The condensed-cyclic compound of claim 1, wherein L₁ and L₂ are each independently selected from groups represented by Formulae 4-11, 4-13, 4-27, and 4-29 to 4-35,

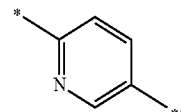
L₁₁ is selected from groups represented by Formulae 4-1, 4-3, 4-5, 4-7 to 4-13, 4-17, and 4-24 to 4-35, and

L₁₂ is selected from groups represented by Formulae 4-1 to 4-35:

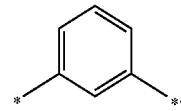
Formula 4-1



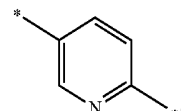
Formula 4-2



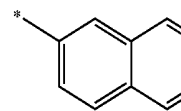
Formula 4-3



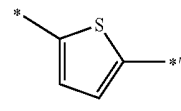
Formula 4-4



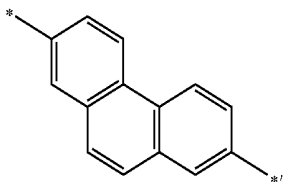
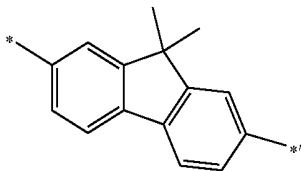
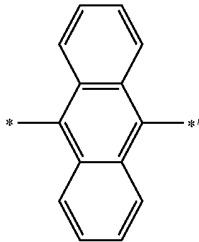
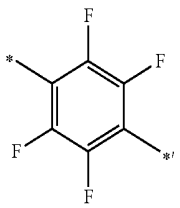
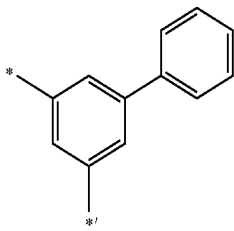
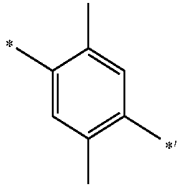
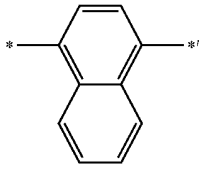
Formula 4-5



Formula 4-6



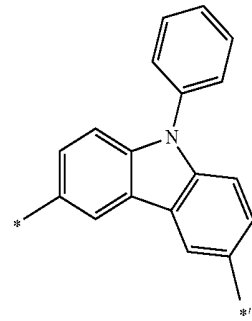
179
-continued



180
-continued

Formula 4-7

5

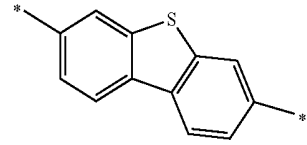


Formula 4-8

10

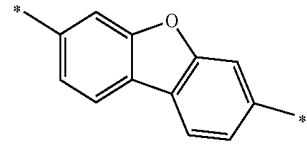
Formula 4-9

15



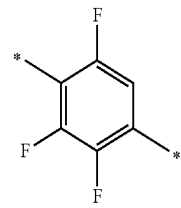
20

25



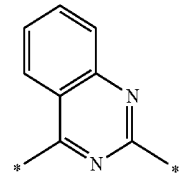
Formula 4-10

30



Formula 4-11

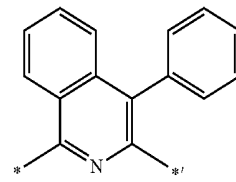
35



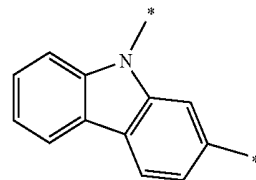
40

Formula 4-12

50

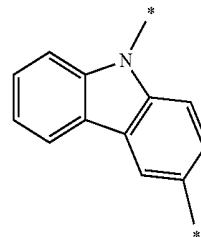


55



Formula 4-13

60



65

Formula 4-14

Formula 4-15

Formula 4-16

Formula 4-17

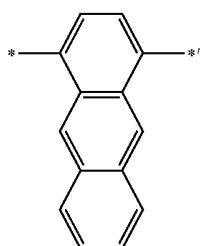
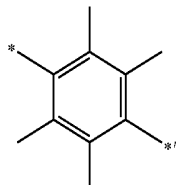
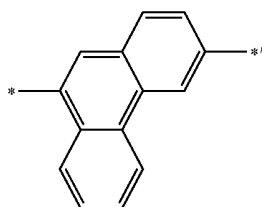
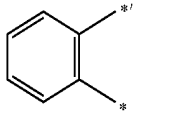
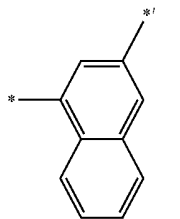
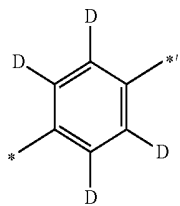
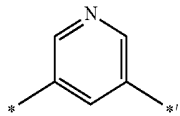
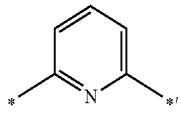
Formula 4-18

Formula 4-19

Formula 4-20

Formula 4-21

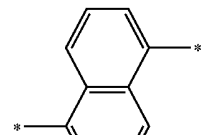
181
-continued



182
-continued

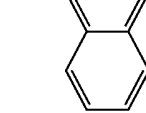
Formula 4-22

5



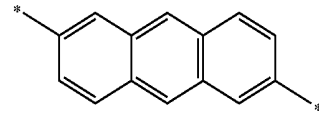
Formula 4-23

10



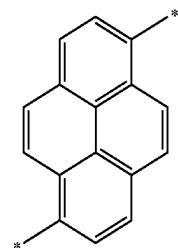
Formula 4-24

15



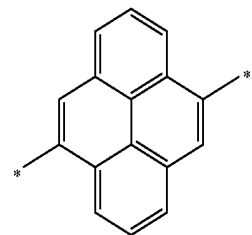
Formula 4-25

20



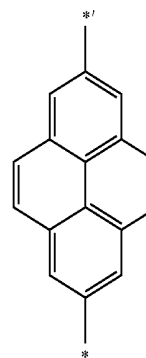
Formula 4-26

35



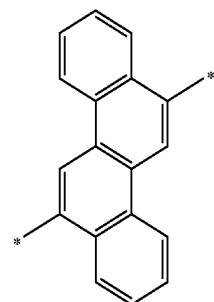
Formula 4-27

40



Formula 4-28

50



Formula 4-29

60

Formula 4-30

Formula 4-31

Formula 4-32

Formula 4-33

Formula 4-34

Formula 4-35

wherein, in Formulae 4-1 to 4-35, * and *¹ each indicate a binding site to an adjacent atom.

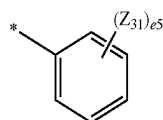
7. The condensed-cyclic compound of claim 1, wherein R₁₁ is selected from

183

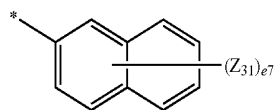
a phenyl group, a biphenyl group, a terphenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenyl group, a heptalenyl group, an indacenyl group, an acenaphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthrenyl 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, and an ovalenyl group;

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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthrenyl 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, and an ovalenyl 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 cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, 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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthrenyl 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, and —Si(Q₃₃)(Q₃₄)(Q₃₅), wherein Q₃₁ to Q₃₃ are each independently 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.

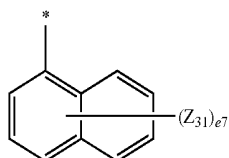
8. The condensed-cyclic compound of claim 1, wherein R₁₁ is selected from groups represented by Formulae 5-1 to 5-19:



Formula 5-1



Formula 5-2

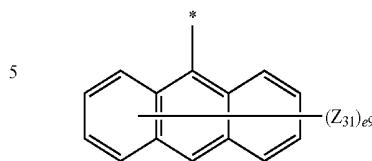


Formula 5-3

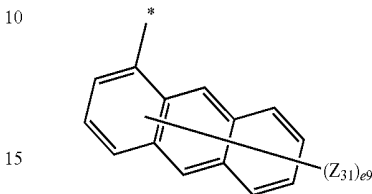
184

-continued

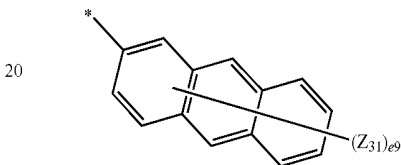
Formula 5-4



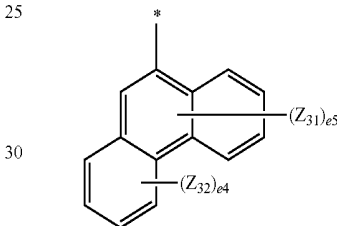
Formula 5-5



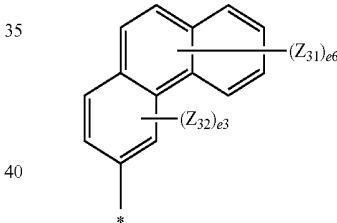
Formula 5-6



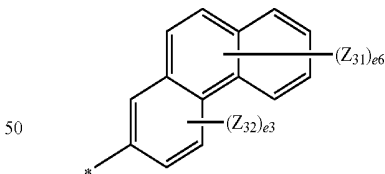
Formula 5-7



Formula 5-8



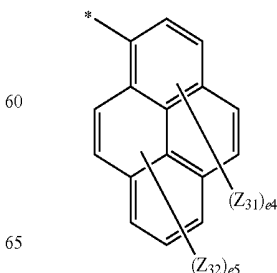
Formula 5-9



50

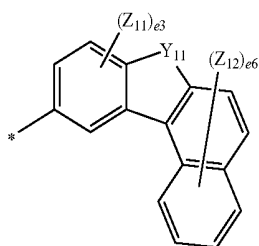
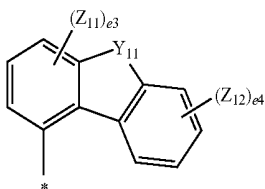
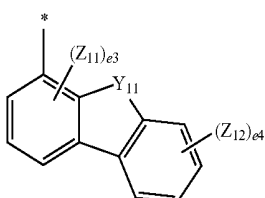
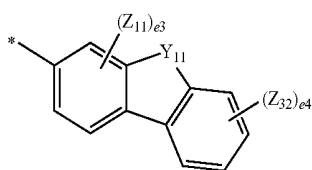
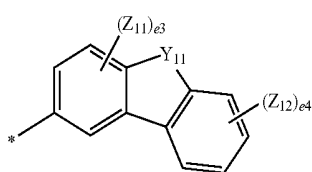
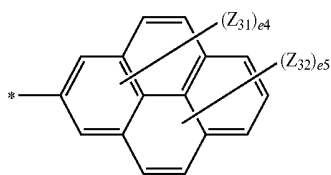
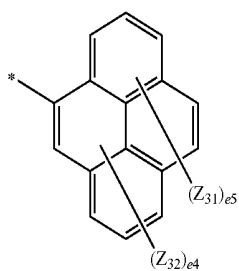
55

Formula 5-10



65

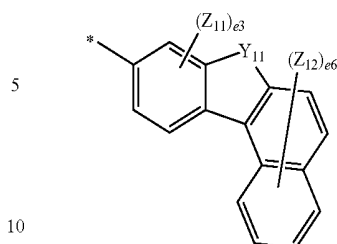
185
-continued



186

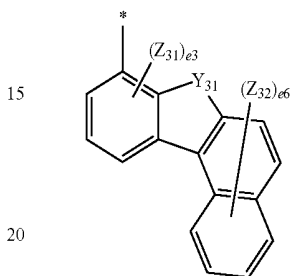
-continued

Formula 5-11



5

Formula 5-12



15

Formula 5-13

20

wherein, in Formulae 5-1 to 5-19,

Y₁₁ is C(Z₁₃)(Z₁₄),

Z₁₁ to Z₁₄ 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 C₁-C₁₀ alkoxy group, a phenyl group, a naphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenanthrenyl group, an anthracenyl group, a pyrenyl group, a chrysenyl group, a biphenyl group, a terphenyl group, and —Si(Q₃₁)(Q₃₂)(Q₃₃),

Formula 5-14

30

Formula 5-15

35

wherein Q₃₁ to Q₃₃ are each independently 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,

40

e₃ is an integer selected from 1 to 3,

e₄ is an integer selected from 1 to 4,

e₅ is an integer selected from 1 to 5,

e₆ is an integer selected from 1 to 6,

e₇ is an integer selected from 1 to 7,

45

e₉ is an integer selected from 1 to 9, and

Formula 5-16

50

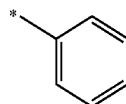
* indicates a binding site to an adjacent atom.

9. The condensed-cyclic compound of claim 1, wherein R₁₁ is selected from groups represented by Formulae 6-1 to 6-27:

50

Formula 6-1

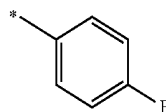
55



Formula 6-2

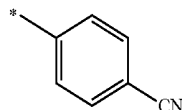
Formula 5-17

60

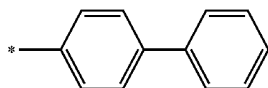
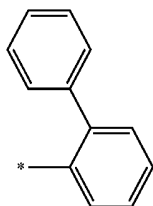
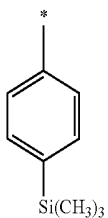
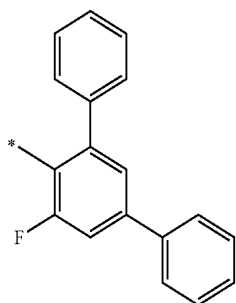
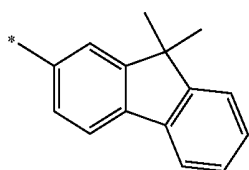
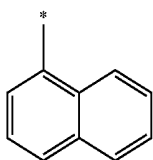
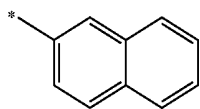
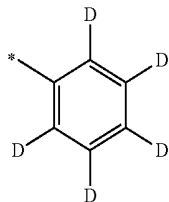
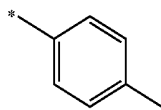


Formula 6-3

65



187
-continued



188
-continued

Formula 6-4

5

Formula 6-5

10

Formula 6-6

15

Formula 6-7

20

Formula 6-8

25

30

Formula 6-9

35

40

Formula 6-10

45

50

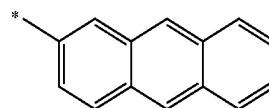
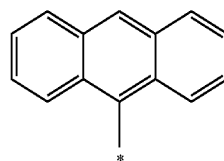
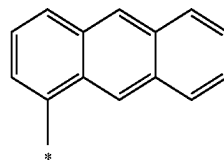
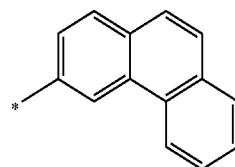
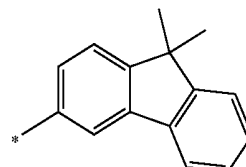
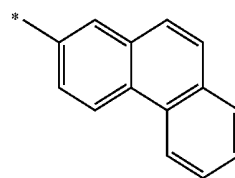
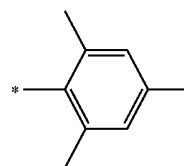
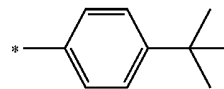
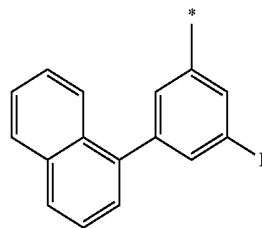
Formula 6-11

55

60

Formula 6-12

65



Formula 6-13

Formula 6-14

Formula 6-15

Formula 6-16

Formula 6-17

Formula 6-18

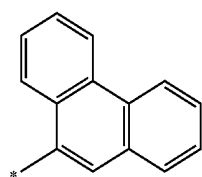
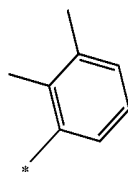
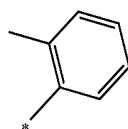
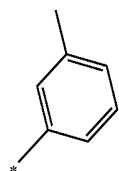
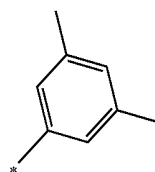
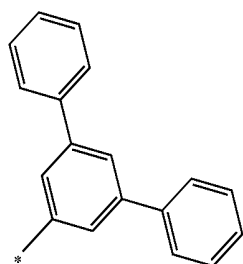
Formula 6-19

Formula 6-20

Formula 6-21

189

-continued



wherein, in Formulae 6-1 to 6-27, * indicates a binding site to an adjacent atom.

10. The condensed-cyclic compound of claim 1, wherein R_1 to R_4 , R_6 and R_{12} to R_{13} 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 C_1 - C_{20} alkyl group, and a C_1 - C_{20} alkoxy group;

a C_1 - C_{20} alkyl group and a C_1 - C_{20} alkoxy group, each substituted with at least one selected from deuterium, —F, —Cl, —Br, —I, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group;

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

190

Formula 6-22

5

10

Formula 6-23

15

Formula 6-24

20

25

Formula 6-25

30

Formula 6-26

35

Formula 6-27

40

45

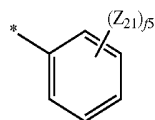
group, an acenaphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexaceny group, a pentaceny group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a 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, a dibenzosilolyl group, a thiadiazolyl group, an imidazopyridinyl group, and an imidazopyrimidinyl group;

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-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexaceny group, a pentaceny group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a 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, a dibenzosilolyl group, a thiadiazolyl group, an imidazopyridinyl group, and an imidazopyrimidinyl 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 cyclopentyl group, a cyclohexyl group, a cycloheptyl group, a cyclopentenyl group, a cyclohexenyl group, a phenyl

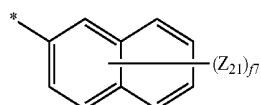
group, a biphenyl group, a terphenyl group, a pentalenyl group, an indenyl group, a naphthyl group, an azulenyl group, a heptalenyl group, an indacenyl group, an acenaphthyl group, a fluorenyl group, a spiro-fluorenyl group, a benzofluorenyl group, a dibenzofluorenyl group, a phenalenyl group, a phenanthrenyl group, an anthracenyl group, a fluoranthenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, a naphthacenyl group, a picenyl group, a perylenyl group, a pentaphenyl group, a hexacenyl group, a pentacenyl group, a rubicenyl group, a coronenyl group, an ovalenyl group, a pyrrolyl group, a thiophenyl group, a furanyl group, an imidazolyl group, a pyrazolyl group, a thiazolyl group, an isothiazolyl group, an oxazolyl group, an isoxazolyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a benzoquinolinyl group, a phthalazinyl group, a naphthyridinyl group, a quinoxalinyl group, a quinazolinyl group, a cinnolinyl group, a carbazolyl group, a phenanthridinyl group, an acridinyl group, a phenanthrolinyl group, a phenazinyl group, a 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, a thiadiazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, and $-\text{Si}(\text{Q}_{31})(\text{Q}_{32})(\text{Q}_{33})$; and $-\text{Si}(\text{Q}_1)(\text{Q}_2)(\text{Q}_3)$.

wherein Q_1 to Q_3 and Q_{31} to Q_{33} are each independently 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.

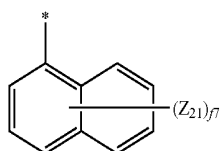
11. The condensed-cyclic compound of claim 1, wherein R_1 to R_4 , R_6 , R_{12} , and R_{13} are each independently selected from hydrogen, deuterium, $-\text{F}$, $-\text{Cl}$, $-\text{Br}$, $-\text{I}$, a hydroxyl group, a cyano group, a nitro group, an amidino group, a hydrazino group, a hydrazono group, a C_1 - C_{10} alkyl group, a C_1 - C_{10} alkoxy group, groups represented by Formulae 7-1 to 7-75, and $-\text{Si}(\text{Q}_1)(\text{Q}_2)(\text{Q}_3)$, wherein Q_1 to Q_3 are each independently 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:



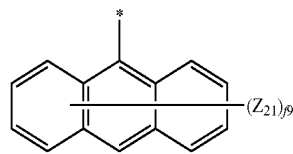
Formula 7-1



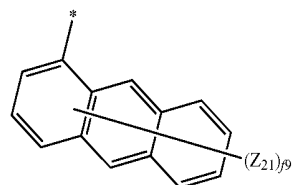
Formula 7-2



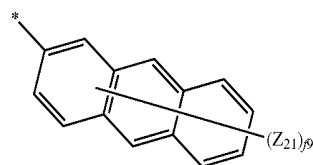
Formula 7-3



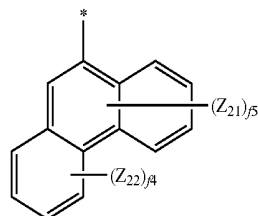
Formula 7-4



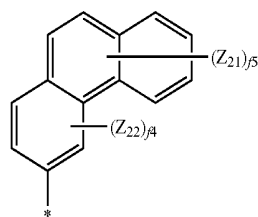
Formula 7-5



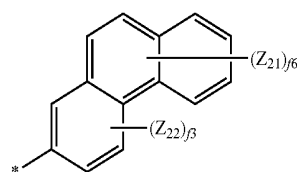
Formula 7-6



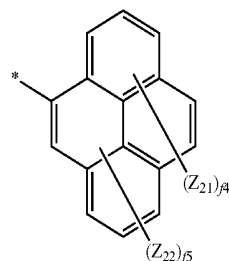
Formula 7-7



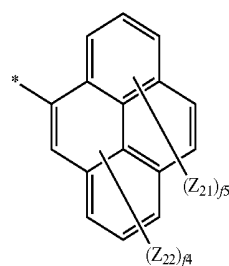
Formula 7-8



Formula 7-9



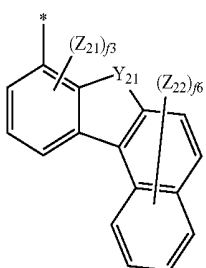
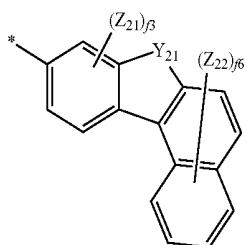
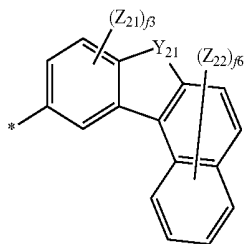
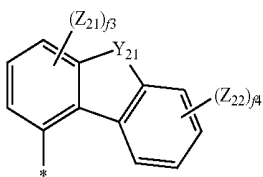
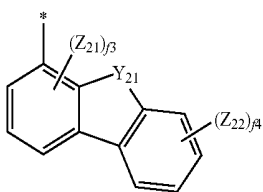
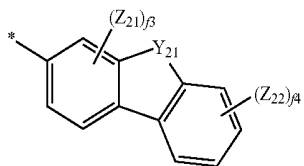
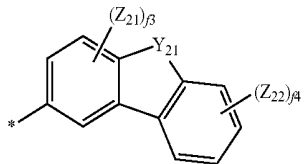
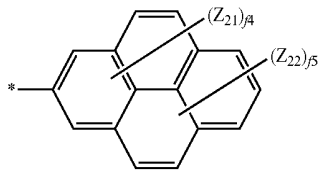
Formula 7-10



Formula 7-11

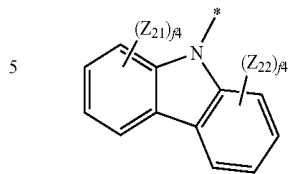
65

193
-continued

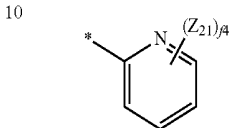


194
-continued

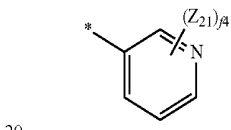
Formula 7-12



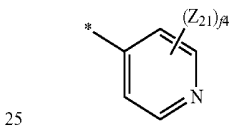
Formula 7-13



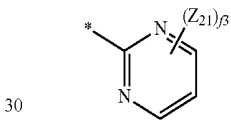
Formula 7-14



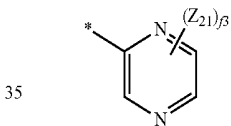
Formula 7-15



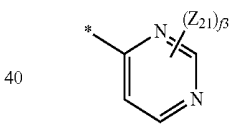
Formula 7-16



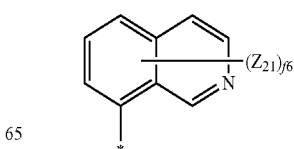
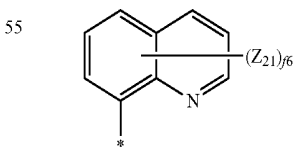
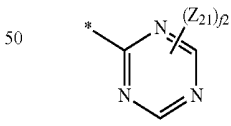
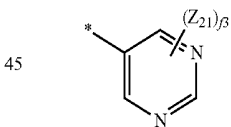
Formula 7-17



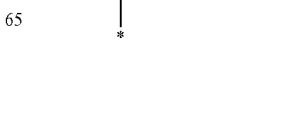
Formula 7-18



Formula 7-19



60



Formula 7-20

Formula 7-21

Formula 7-22

Formula 7-23

Formula 7-24

Formula 7-25

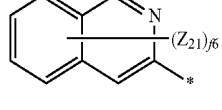
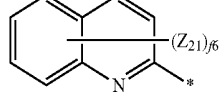
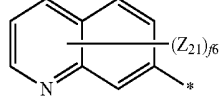
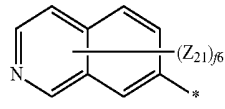
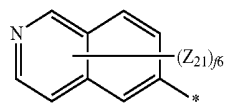
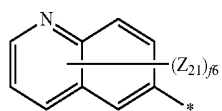
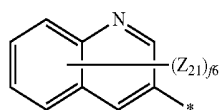
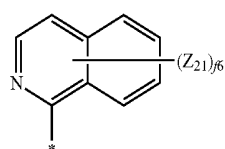
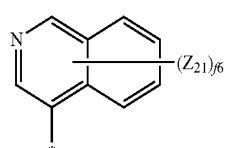
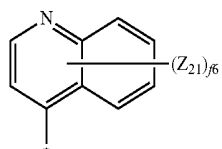
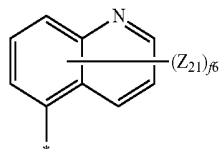
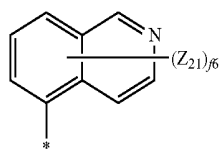
Formula 7-26

Formula 7-27

Formula 7-28

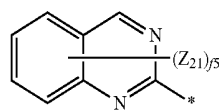
Formula 7-29

Formula 7-30



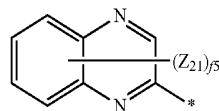
Formula 7-31

5



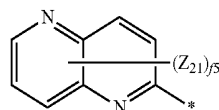
Formula 7-32

10



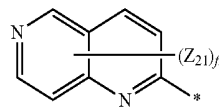
Formula 7-33

15



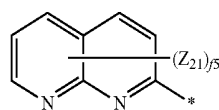
Formula 7-34

25



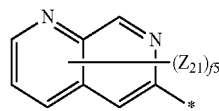
Formula 7-35

30



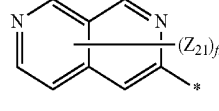
Formula 7-36

35



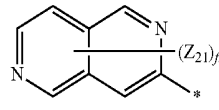
Formula 7-37

40



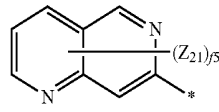
Formula 7-38

45



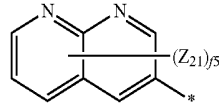
Formula 7-39

50



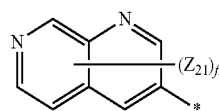
Formula 7-40

55



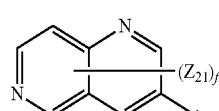
Formula 7-41

60



Formula 7-42

65



Formula 7-43

Formula 7-44

Formula 7-45

Formula 7-47

Formula 7-48

Formula 7-49

Formula 7-50

Formula 7-51

Formula 7-52

Formula 7-53

Formula 7-54

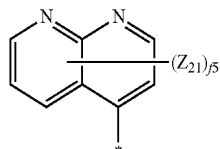
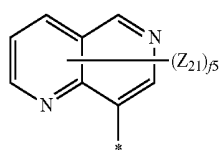
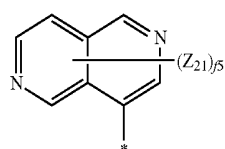
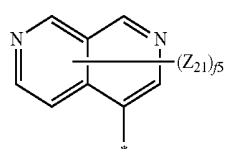
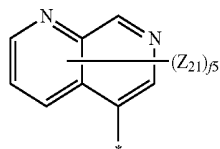
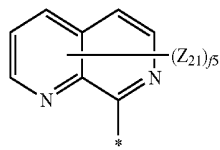
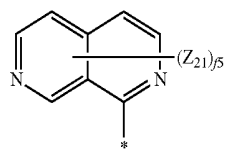
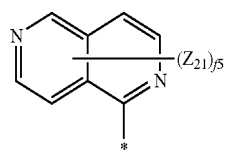
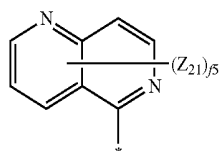
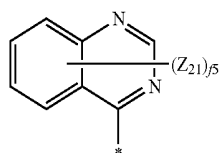
Formula 7-55

Formula 7-56

Formula 7-57

197

-continued

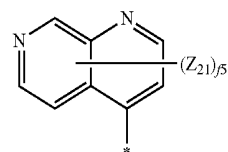


198

-continued

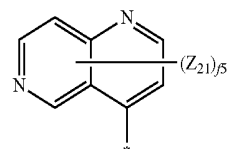
Formula 7-58

5



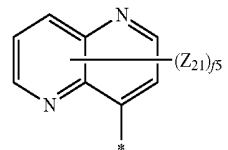
Formula 7-59

10



Formula 7-60

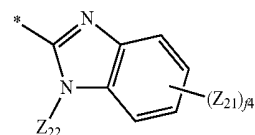
15



20

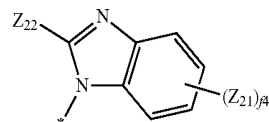
Formula 7-61

25



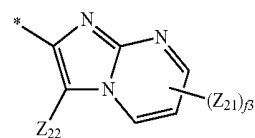
Formula 7-62

30



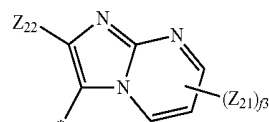
Formula 7-62

35



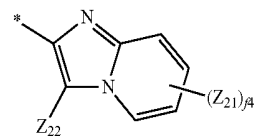
Formula 7-63

40



Formula 7-64

50



Formula 7-65

55

wherein, in Formulae 7-1 to 7-75,
 Y_{21} is selected from O, S, $C(Z_{23})(Z_{24})$, $N(Z_{25})$, and
 $Si(Z_{26})(Z_{27})$,

Formula 7-66

60

Z_{21} to Z_{27} 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 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-fluorenyl group, a benzofluorenyl
 group, a dibenzofluorenyl group, a phenanthrenyl
 group, an anthracenyl group, a pyrenyl group, a chrysenyl
 group, a pyridinyl group, a pyrazinyl group, a

65

Formula 7-67

Formula 7-68

Formula 7-69

Formula 7-70

Formula 7-71

Formula 7-72

Formula 7-73

Formula 7-74

Formula 7-75

199

pyrimidinyl group, a pyridazinyl group, a quinolinyl group, an isoquinolinyl group, a quinoxalinyl group, a quinazolinyl group, a carbazolyl group, a triazinyl group, and $-\text{Si}(\text{Q}_{31})(\text{Q}_{32})(\text{Q}_{33})$, wherein Q_{31} to Q_{33} are each independently selected from a 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,

f2 is an integer selected from 1 and 2,

f3 is an integer from 1 to 3,

f4 is an integer from 1 to 4,

f5 is an integer from 1 to 5,

f6 is an integer from 1 to 6,

f7 is an integer from 1 to 7,

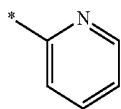
f8 is an integer from 1 to 8,

f9 is an integer from 1 to 9, and

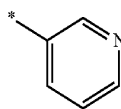
* indicates a binding site to an adjacent atom.

12. The condensed-cyclic compound of claim 1, wherein R_3 , R_4 , R_6 and R_{13} are each independently selected from hydrogen, 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_{10} alkyl group, a C_1 - C_{10} alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, a naphthyl group, and $-\text{Si}(\text{Q}_1)(\text{Q}_2)(\text{Q}_3)$, and

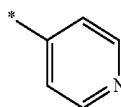
R_1 , R_2 , and R_{12} are each independently selected from groups represented by Formulae 8-1 to 8-159, wherein Q_1 to Q_3 are each independently selected from a C_1 - C_{10} alkyl group, a C_1 - C_{10} alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group:



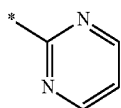
Formula 8-1



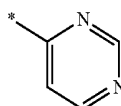
Formula 8-2



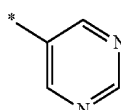
Formula 8-3



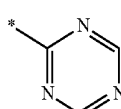
Formula 8-4



Formula 8-5



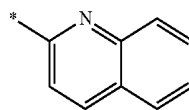
Formula 8-6



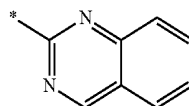
Formula 8-7

200

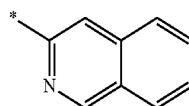
-continued



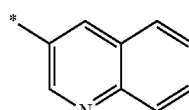
Formula 8-8



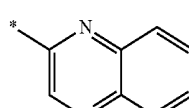
Formula 8-9



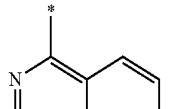
Formula 8-10



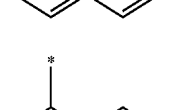
Formula 8-11



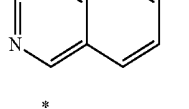
Formula 8-12



Formula 8-13

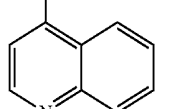


Formula 8-14

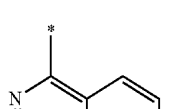


35

Formula 8-15

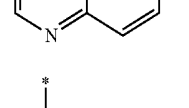


40



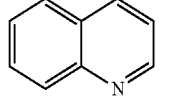
45

Formula 8-16

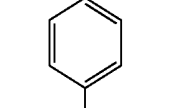


50

Formula 8-17

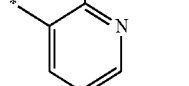


55



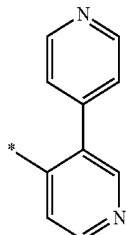
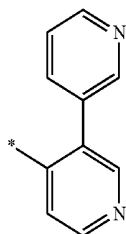
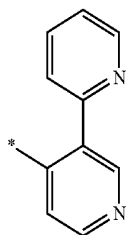
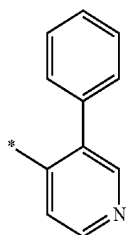
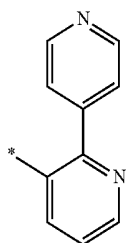
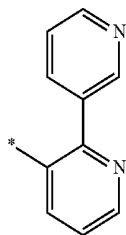
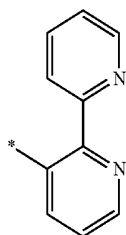
60

Formula 8-18



65

201
-continued



Formula 8-19

5

Formula 8-20

10

Formula 8-21

20

Formula 8-22

30

Formula 8-23

40

Formula 8-24

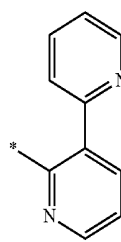
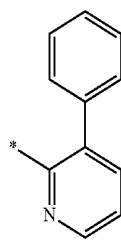
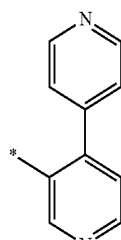
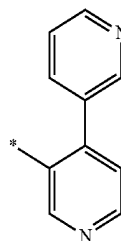
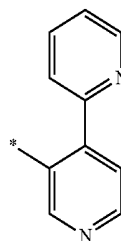
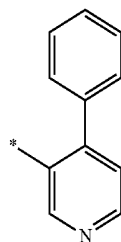
50

Formula 8-25

60

65

202
-continued



Formula 8-26

Formula 8-27

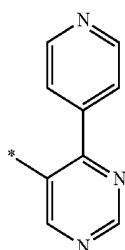
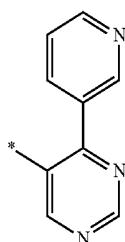
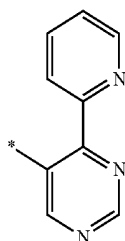
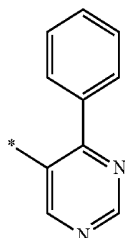
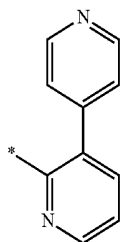
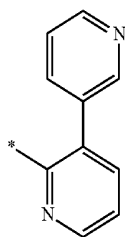
Formula 8-28

Formula 8-29

Formula 8-30

Formula 8-31

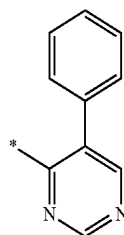
203
-continued



204
-continued

Formula 8-32

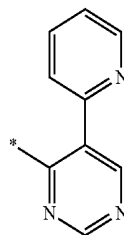
5



10

Formula 8-33

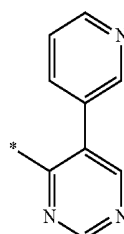
15



20

Formula 8-34

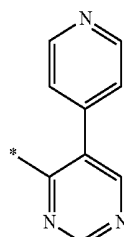
25



30

Formula 8-35

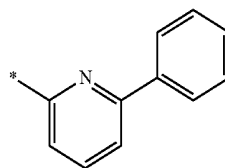
35



40

Formula 8-36

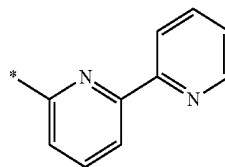
45



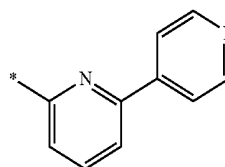
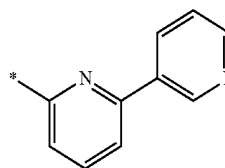
50

Formula 8-37

60



65



Formula 8-38

Formula 8-39

Formula 8-40

Formula 8-41

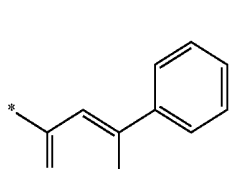
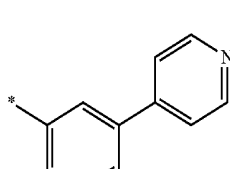
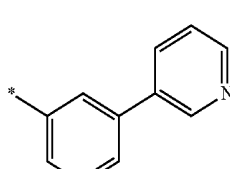
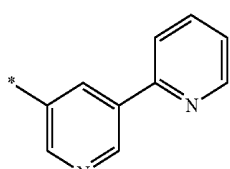
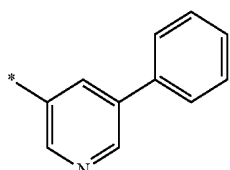
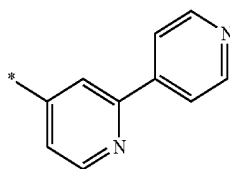
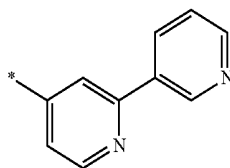
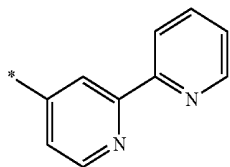
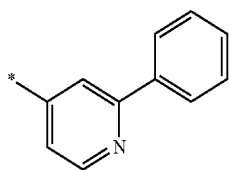
Formula 8-42

Formula 8-43

Formula 8-44

Formula 8-45

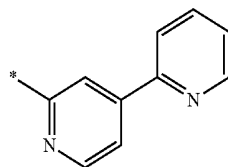
205
-continued



206
-continued

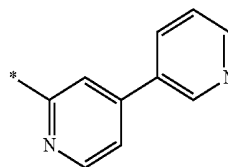
Formula 8-46

5



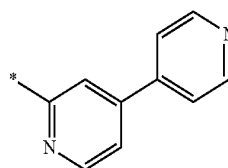
Formula 8-47

10



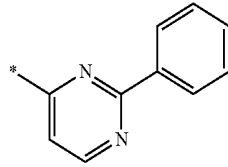
Formula 8-48

15



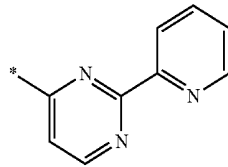
Formula 8-49

20



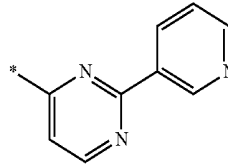
Formula 8-50

25



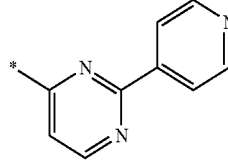
Formula 8-51

30



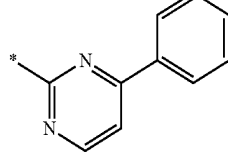
Formula 8-52

35



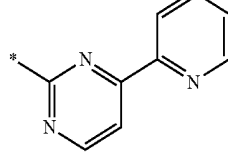
Formula 8-53

40



Formula 8-54

45



Formula 8-55

Formula 8-56

Formula 8-57

Formula 8-58

Formula 8-59

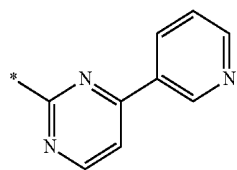
Formula 8-60

Formula 8-61

Formula 8-62

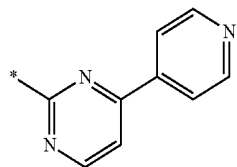
Formula 8-63

207
-continued



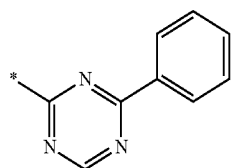
Formula 8-64

5



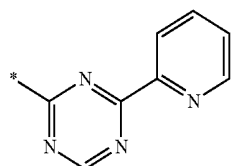
Formula 8-65

10



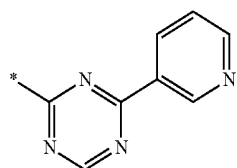
Formula 8-66

15



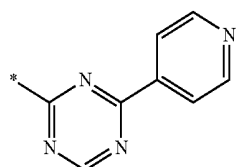
Formula 8-67

25



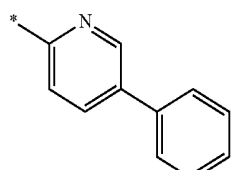
Formula 8-68

30



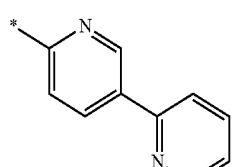
Formula 8-69

40



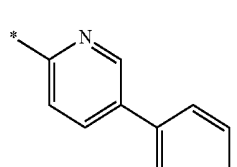
Formula 8-70

45



Formula 8-71

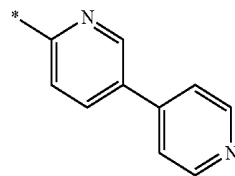
55



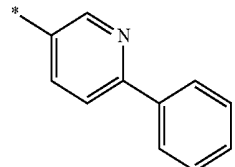
Formula 8-72

60

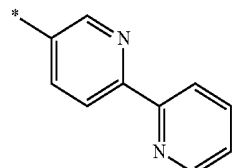
208
-continued



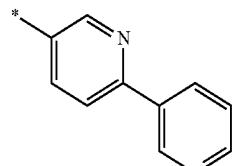
Formula 8-73



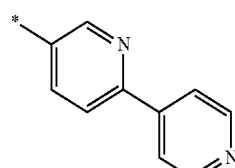
Formula 8-74



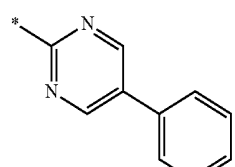
Formula 8-75



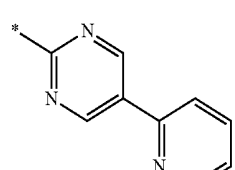
Formula 8-76



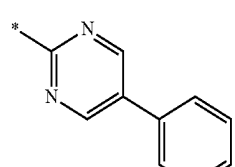
Formula 8-77



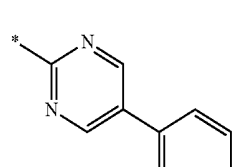
Formula 8-78



Formula 8-79



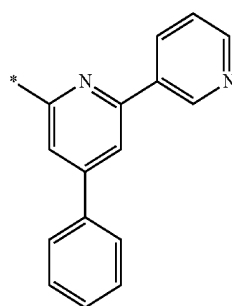
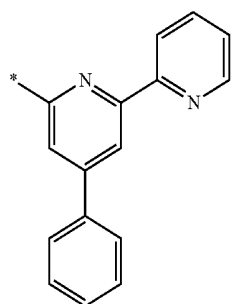
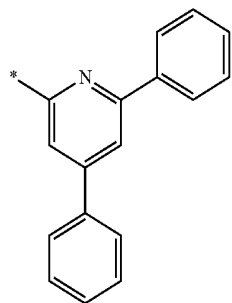
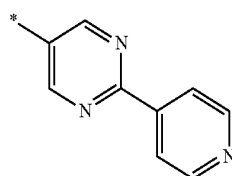
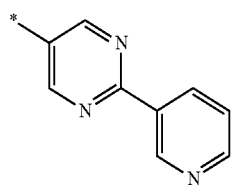
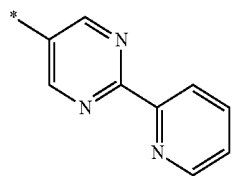
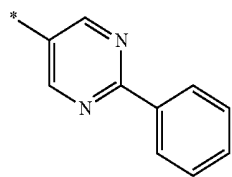
Formula 8-80



Formula 8-81

65

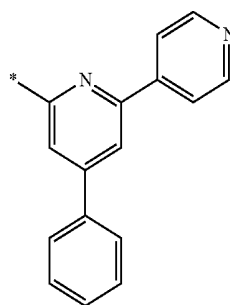
209
-continued



210
-continued

Formula 8-82

5

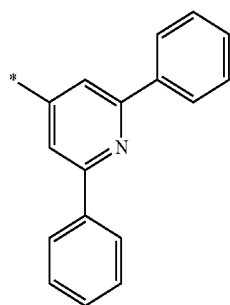


Formula 8-83

10

Formula 8-84

15

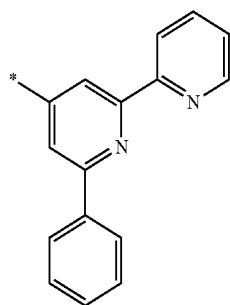


Formula 8-85

25

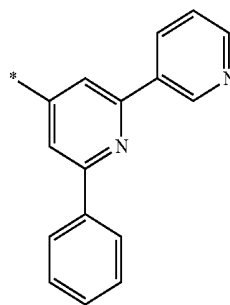
Formula 8-86

30



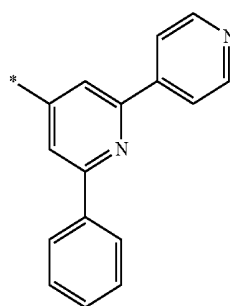
Formula 8-87

45



Formula 8-88

55



Formula 8-89

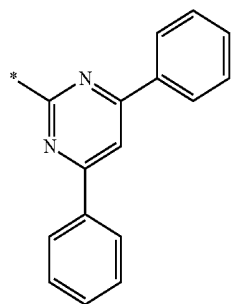
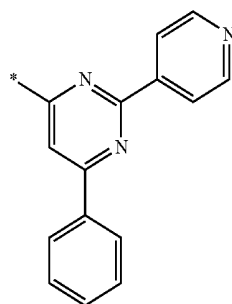
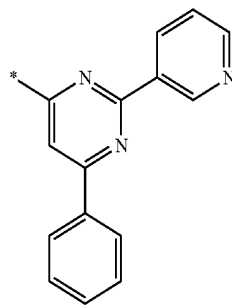
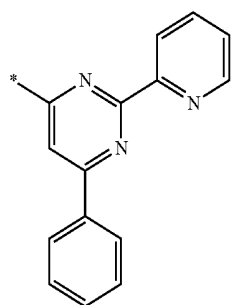
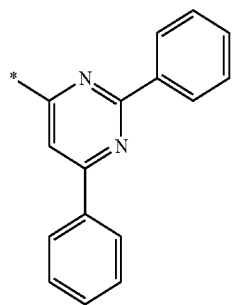
Formula 8-90

Formula 8-91

Formula 8-92

Formula 8-93

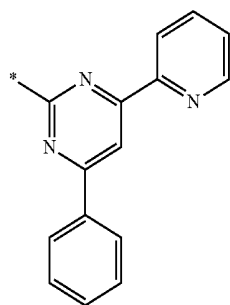
211
-continued



212
-continued

Formula 8-94

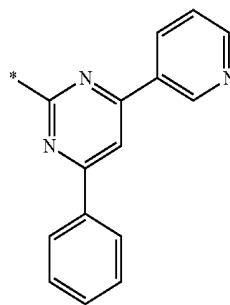
5



10

Formula 8-95

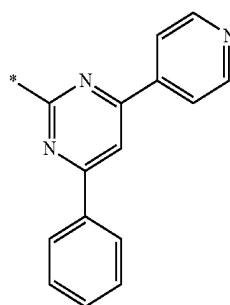
15



20

Formula 8-96

25

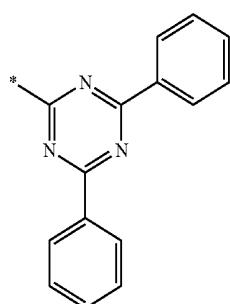


30

35

Formula 8-97

40

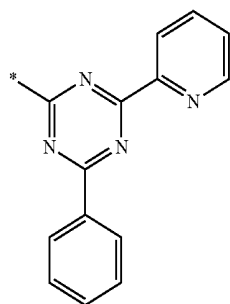


45

50

Formula 8-98

55



60

65

Formula 8-99

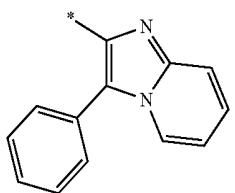
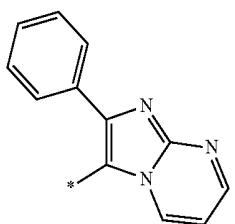
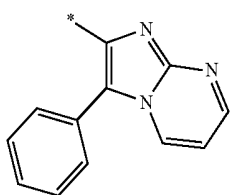
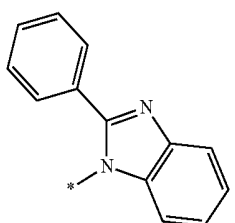
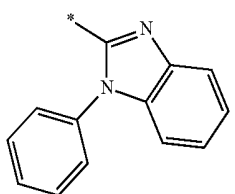
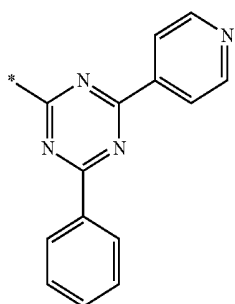
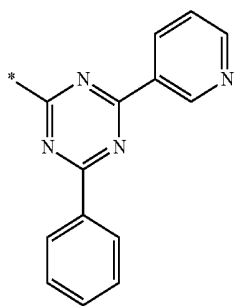
Formula 8-100

Formula 8-101

Formula 8-102

Formula 8-103

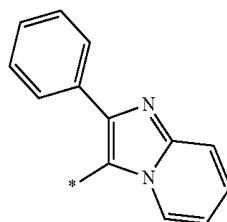
213
-continued



214
-continued

Formula 8-104

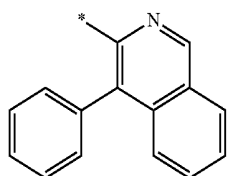
5



10

Formula 8-105

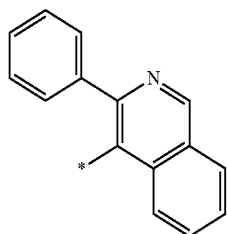
15



20

Formula 8-106

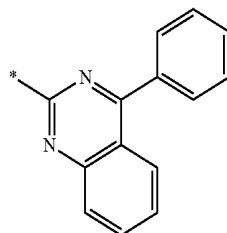
25



30

Formula 8-107

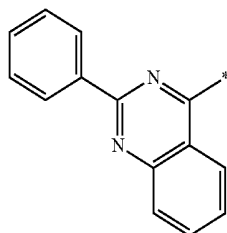
35



40

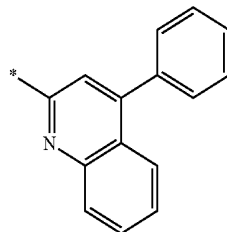
Formula 8-108

45



Formula 8-109

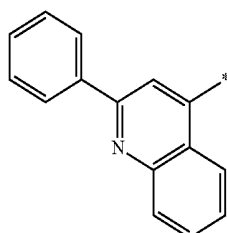
50



55

Formula 8-110

60



65

Formula 8-111

Formula 8-112

Formula 8-113

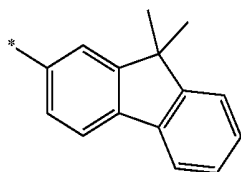
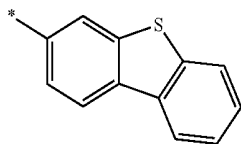
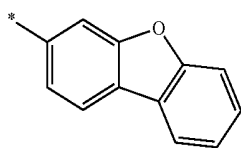
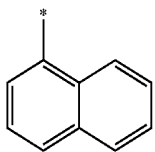
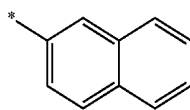
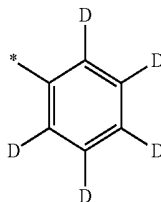
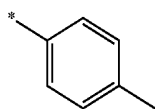
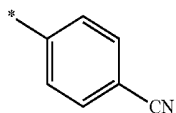
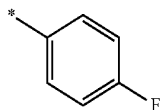
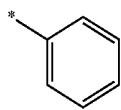
Formula 8-114

Formula 8-115

Formula 8-116

Formula 8-117

215
-continued



216
-continued

Formula 8-118

5

Formula 8-119

10

Formula 8-120

15

Formula 8-121

20

Formula 8-122

25

Formula 8-123

30

Formula 8-124

35

Formula 8-125

45

Formula 8-126

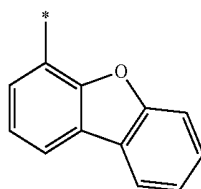
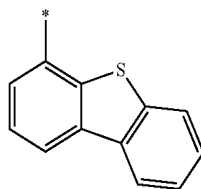
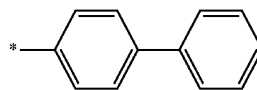
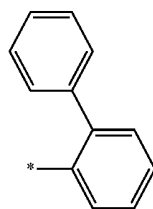
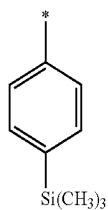
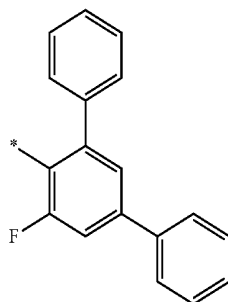
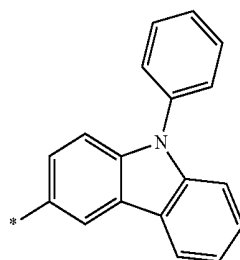
50

55

Formula 8-127

60

65



Formula 8-128

Formula 8-129

Formula 8-130

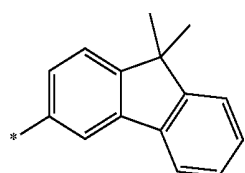
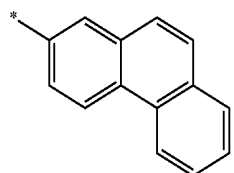
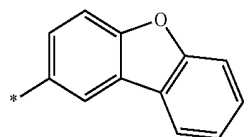
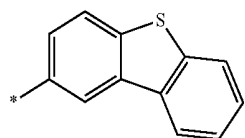
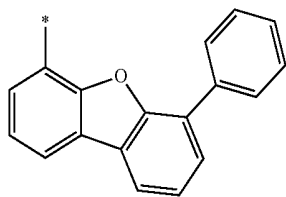
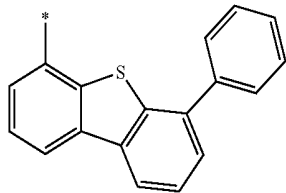
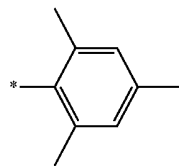
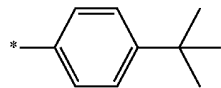
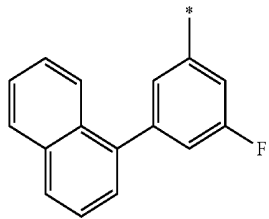
Formula 8-131

Formula 8-132

Formula 8-133

Formula 8-134

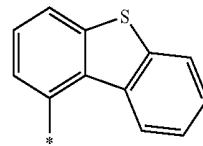
217
-continued



218
-continued

Formula 8-135

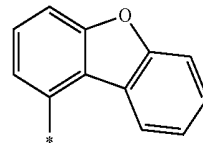
5



Formula 8-144

Formula 8-136

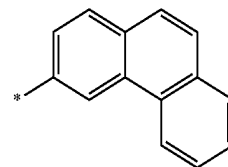
10



Formula 8-145

Formula 8-137

15

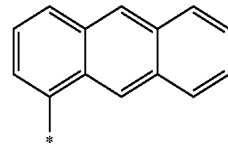


Formula 8-146

20

Formula 8-138

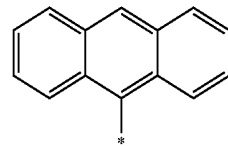
25



Formula 8-147

Formula 8-139

30

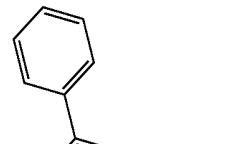


Formula 8-148

35

Formula 8-140

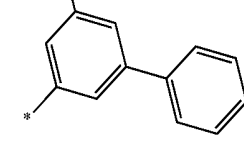
40



Formula 8-149

Formula 8-141

45

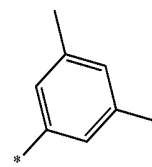


Formula 8-150

50

Formula 8-142

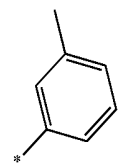
55



Formula 8-151

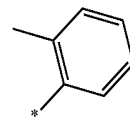
Formula 8-143

60



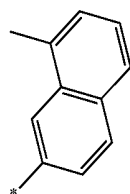
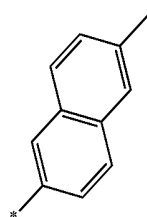
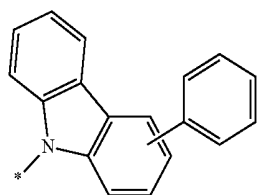
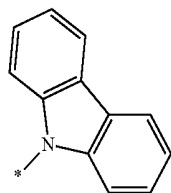
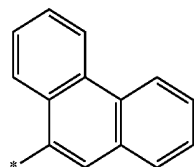
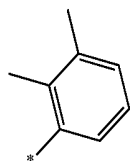
Formula 8-152

65



Formula 8-153

219
-continued



wherein, in Formulae 8-1 to 8-159, * indicates a binding site to an adjacent atom.

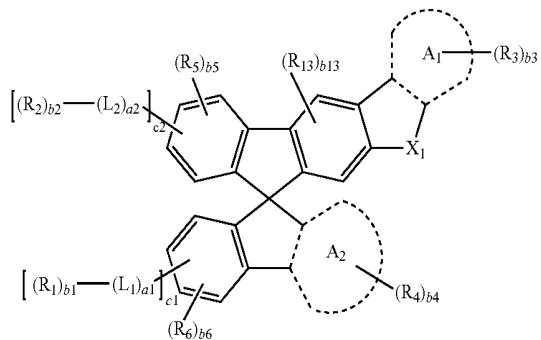
13. The condensed-cyclic compound of claim 1, wherein the condensed-cyclic compound is represented by one of Formulae 1A to 1E:

220

Formula 8-154

5

Formula 1A



Formula 8-155

10

15

Formula 1B

Formula 8-156

20

25

Formula 8-157

30

35

Formula 1C

Formula 8-158

40

45

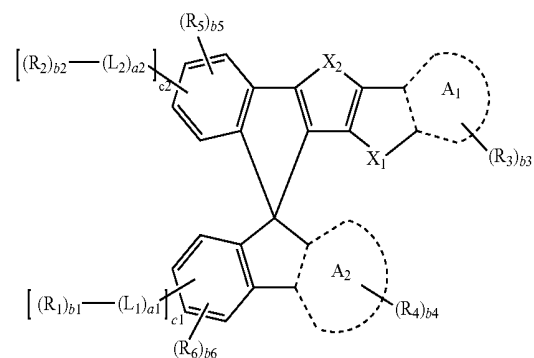
50

Formula 8-159

55

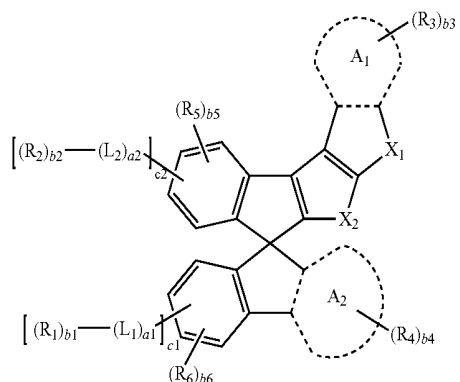
60

Formula 1D



221
-continued

Formula 1E



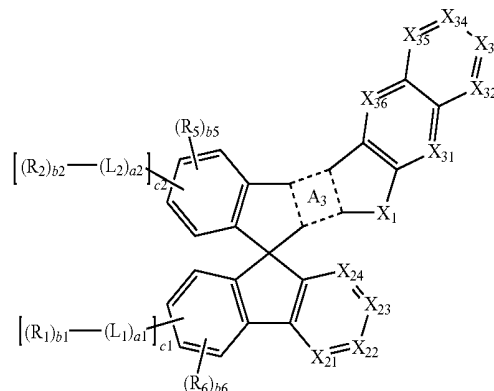
5

10

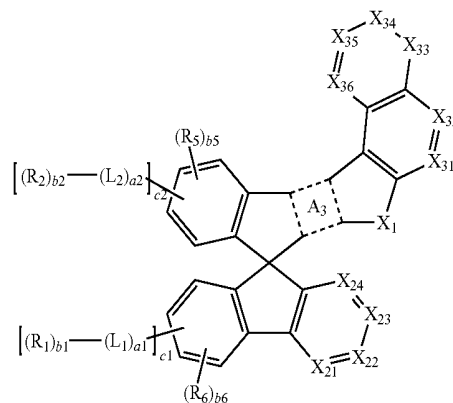
15

222
-continued

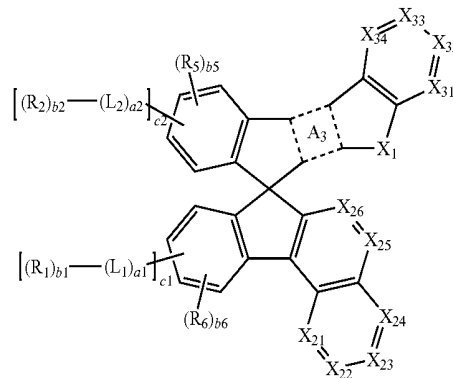
Formula 1-3



Formula 1-4



Formula 1-5



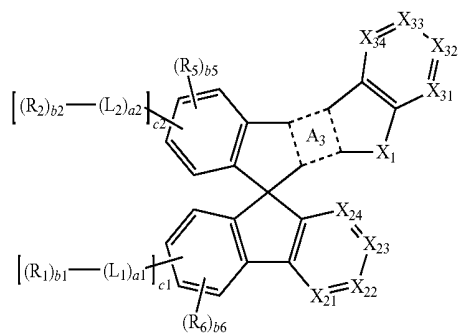
wherein, in Formulae 1A to 1E, ring A₁, ring A₂, X₁, X₂,
L₁, L₂, a₁, a₂, L₁₁, L₁₂, a₁₁, a₁₂, R₁ to R₆, R₁₁ to R₁₃,
b₁ to b₆, b₁₁ to b₁₃, c₁, and c₂ are the same as those
defined in claim 1.

20

14. The condensed-cyclic compound of claim 1, wherein
the condensed-cyclic compound is represented by one of
Formulae 1-1 to 1-7:

25

Formula 1-1



30

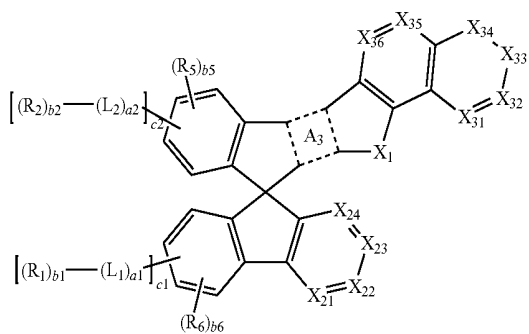
35

40

45

50

Formula 1-2

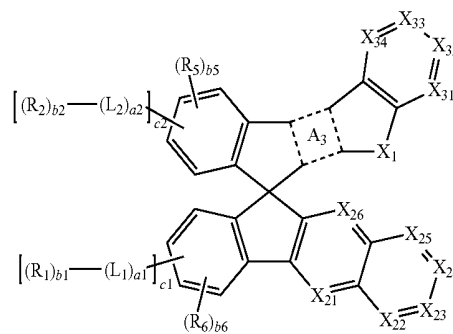


55

60

65

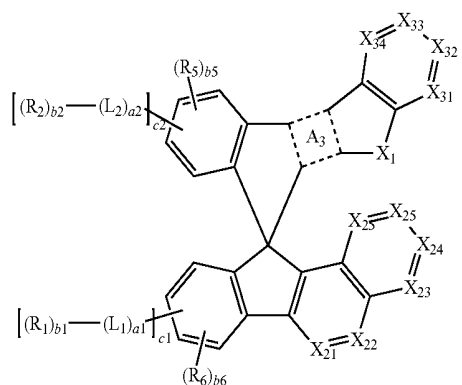
Formula 1-6



223

-continued

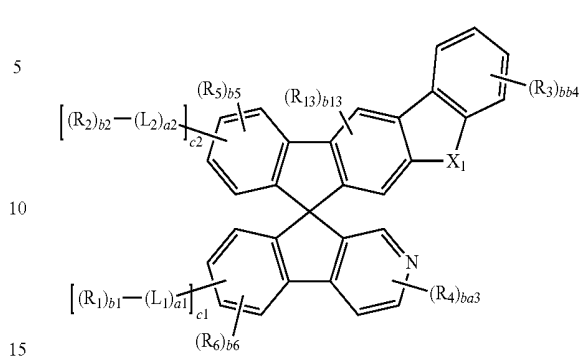
Formula 1-7



224

-continued

Formula 1(3)



Formula 1(4)

wherein, in Formulae 1-1 to 1-7,

ring A₃, X₁, X₂, L₁, L₂, a₁, a₂, L₁₁, L₁₂, a₁₁, a₁₂, R₁, R₂, R₅, R₆, R₁₁ to R₁₃, b₁, b₂, b₅, b₆, b₁₁ to b₁₃, c₁, and c₂ are the same as those defined in claim 1,

X₂₁ is N or C(R₂₁), X₂₂ is N or C(R₂₂), X₂₃ is N or C(R₂₃), X₂₄ is N or C(R₂₄), X₂₅ is N or C(R₂₅), X₂₆ is N or C(R₂₆),

at least one selected from X₂₁ to X₂₆ is N,

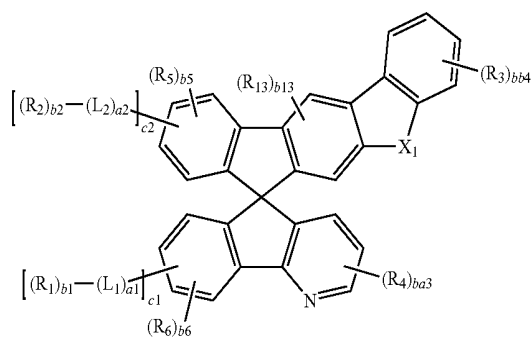
X₃₁ is N or C(R₃₁), X₃₂ is N or C(R₃₂), X₃₃ is N or C(R₃₃), X₃₄ is N or C(R₃₄), X₃₅ is N or C(R₃₅), X₃₆ is N or C(R₃₆),

R₂₁ and R₂₆ are each independently the same as defined in claim 1 with reference to R₃, and

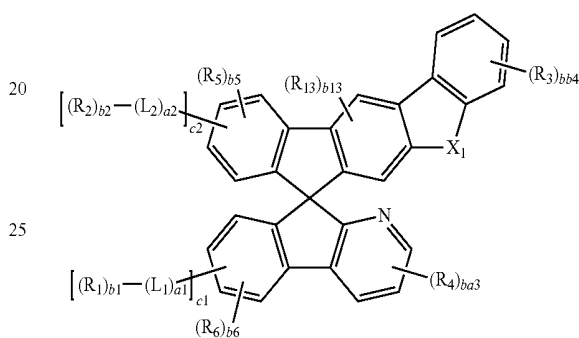
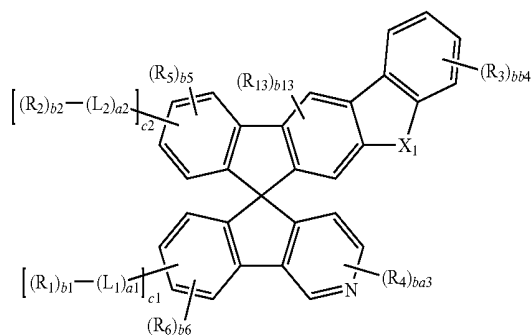
R₃₁ and R₃₆ are each independently the same as defined in claim 1 with reference to R₄.

15. The condensed-cyclic compound of claim 1, wherein the condensed-cyclic compound is represented by one of Formulae 1(1) to 1(19):

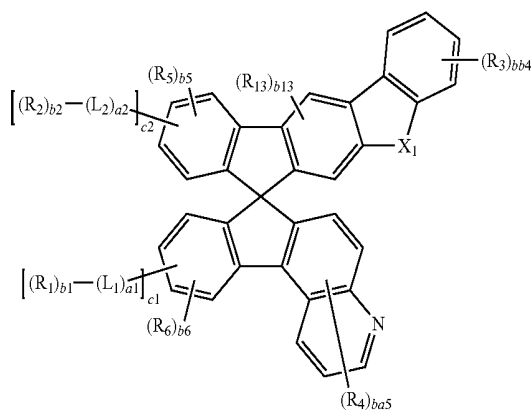
Formula 1(1)



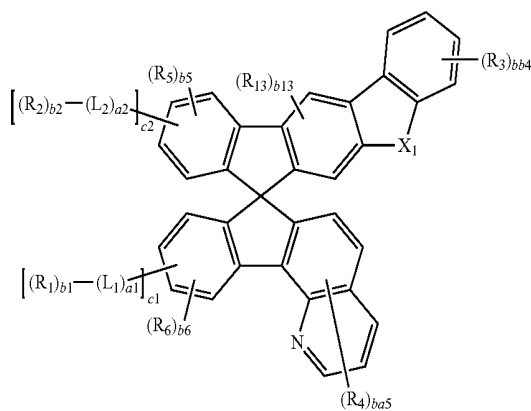
Formula 1(2)



Formula 1(5)

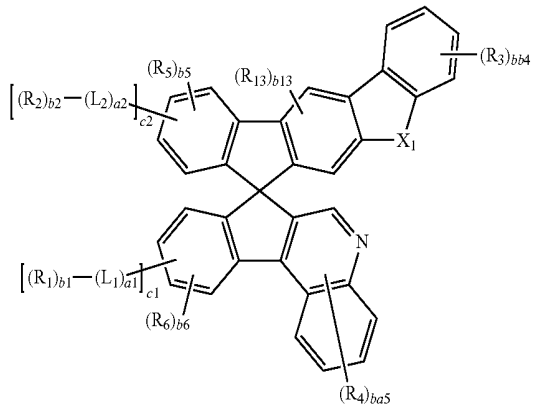


Formula 1(6)



225
-continued

Formula 1(7)



5

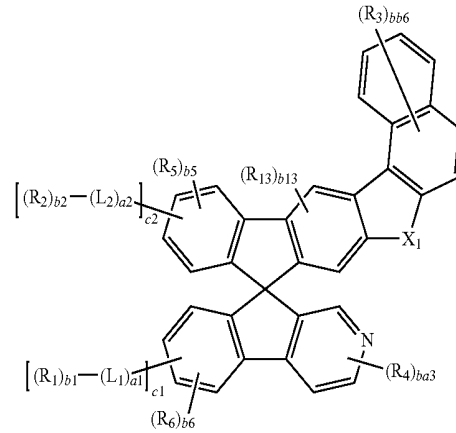
10

15

20

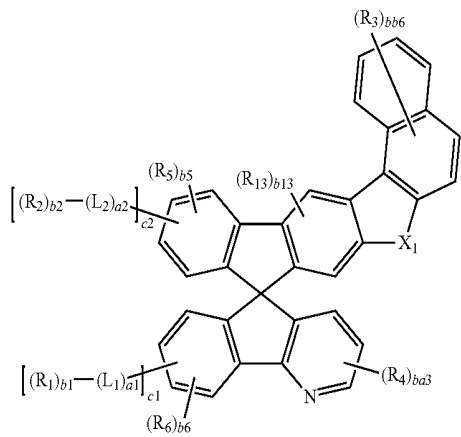
226
-continued

Formula 1(10)



20

Formula 1(8)



25

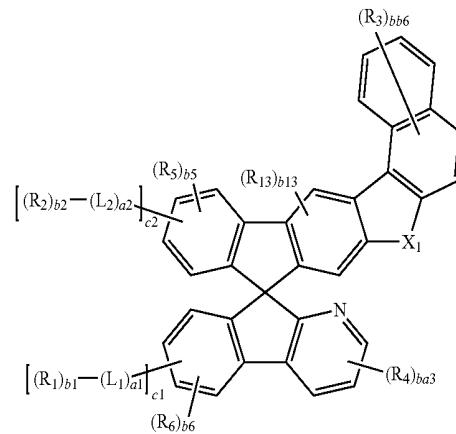
30

35

40

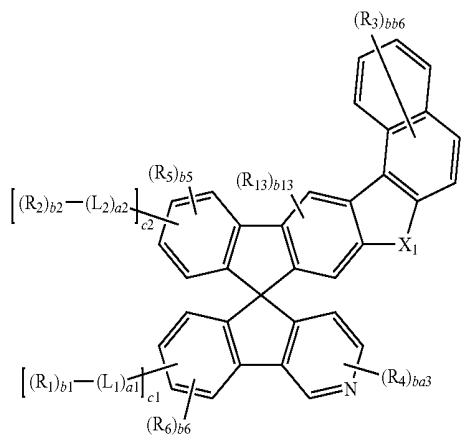
45

Formula 1(11)



45

Formula 1(9)



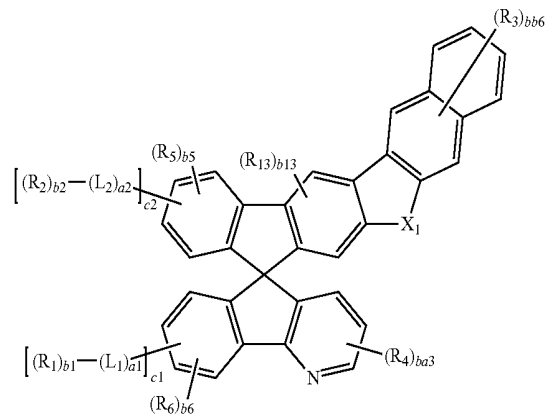
50

55

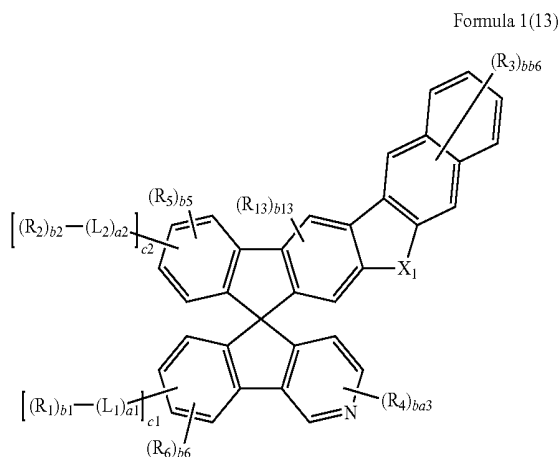
60

65

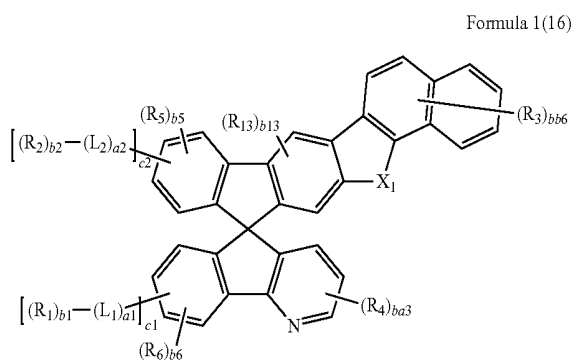
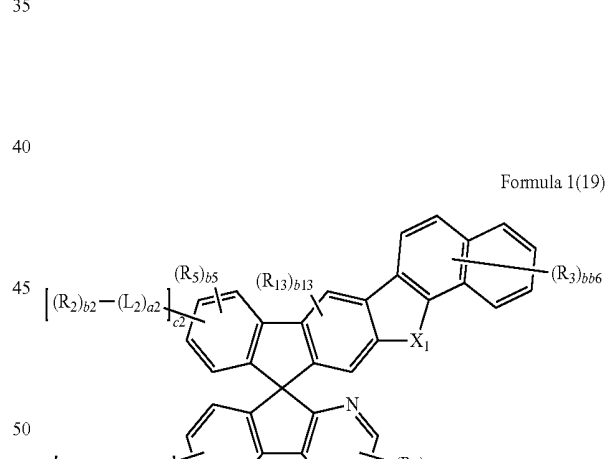
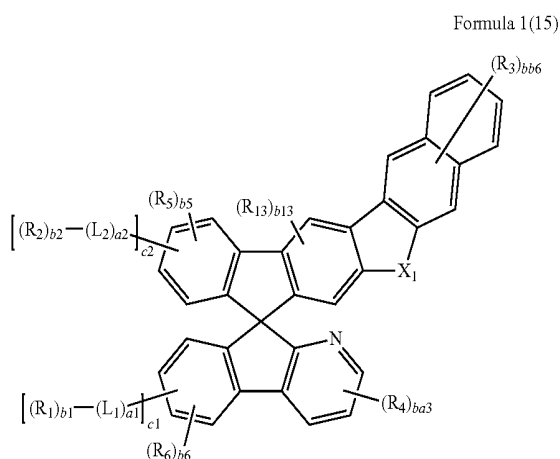
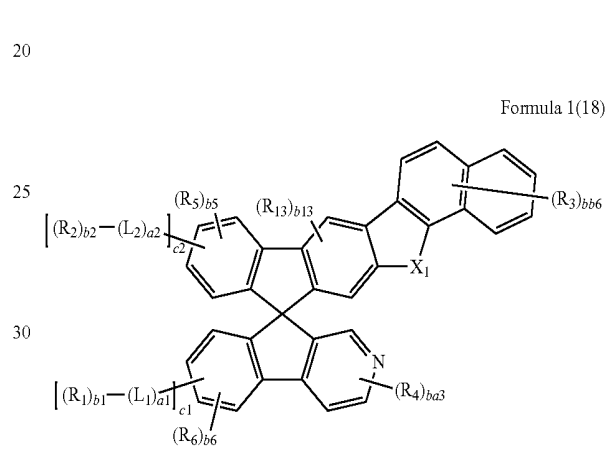
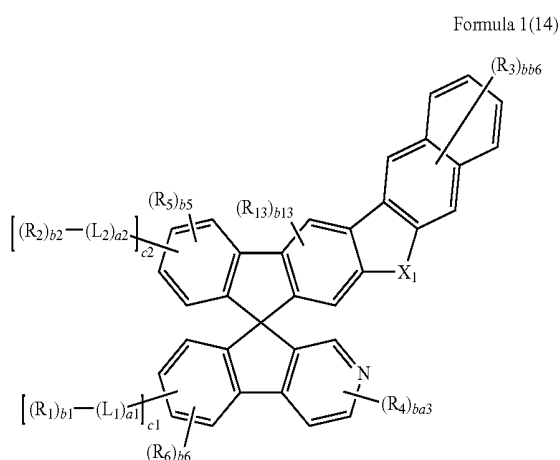
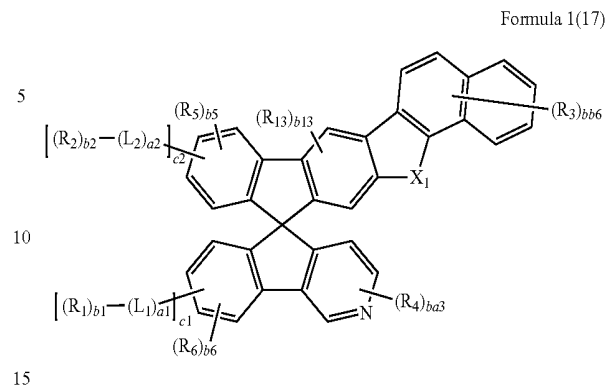
Formula 1(12)



227
-continued



228
-continued



wherein, in Formulae 1(1) to 1(19),

X_1 , L_1 , L_2 , a_1 , a_2 , L_{11} , L_{12} , a_{11} , a_{12} , R_1 to R_6 , R_{11} to R_{13} , b_1 to b_6 , b_{11} to b_{13} , c_1 , and c_2 are the same as those defined in claim 1,

ba_3 is an integer from 0 to 3,

bb_4 is an integer from 0 to 4,

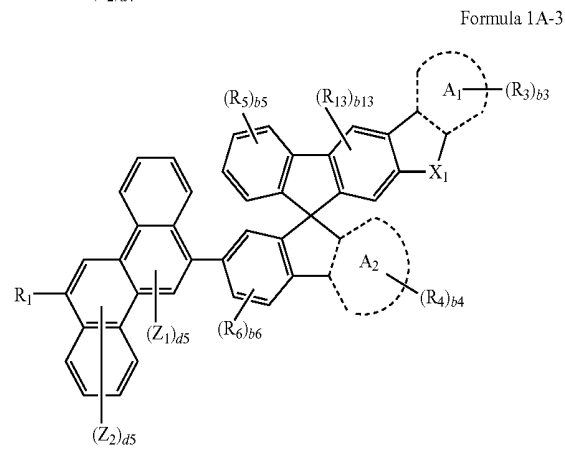
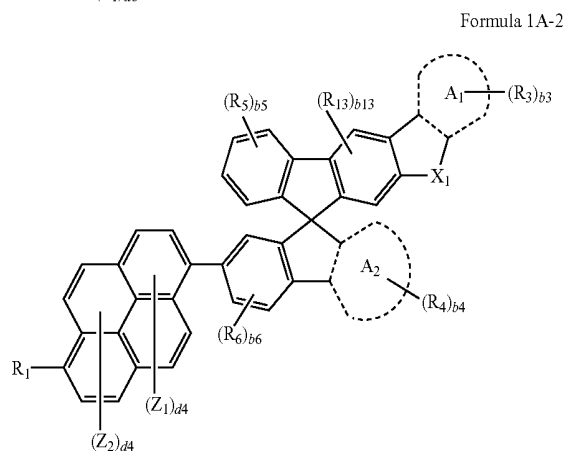
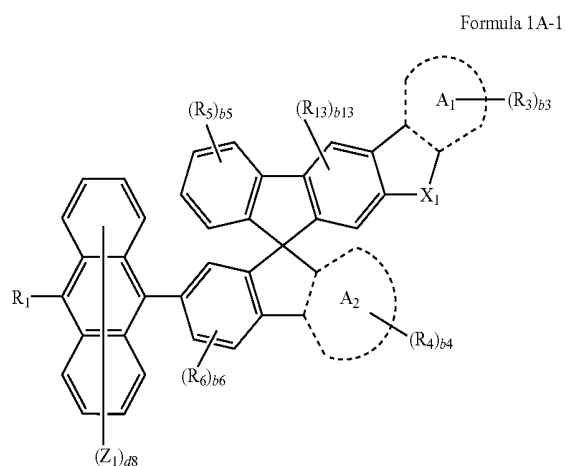
ba_5 is an integer from 0 to 5, and

bb_6 is an integer from 0 to 6.

16. A condensed-cyclic compound represented by one of Formulae 1A-1 to 1A-3:

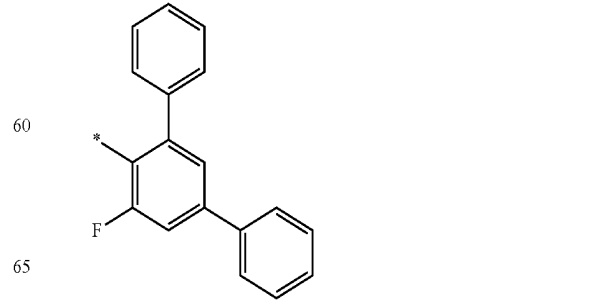
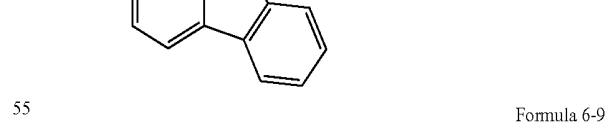
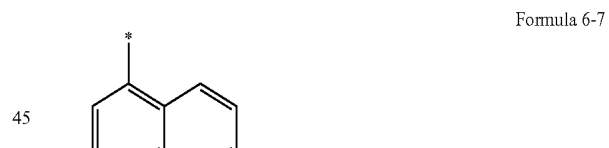
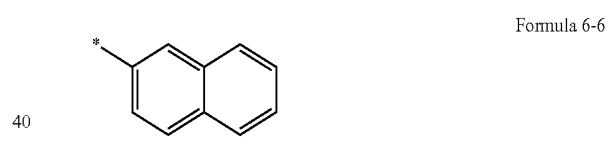
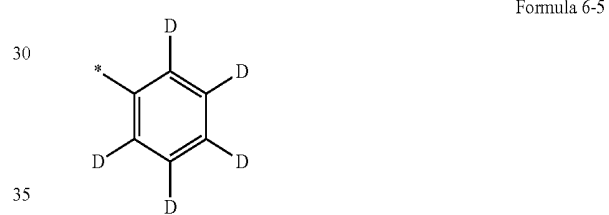
229

230

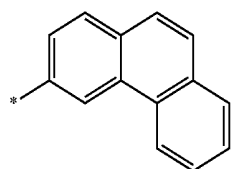
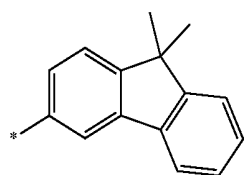
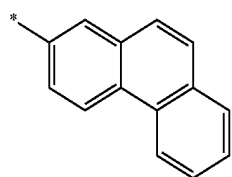
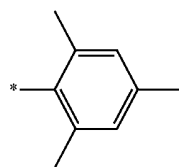
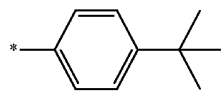
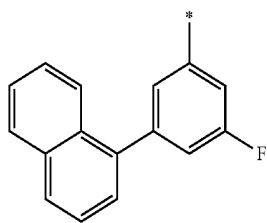
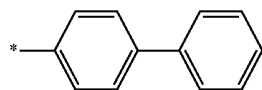
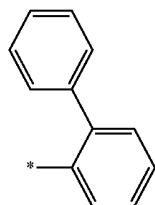
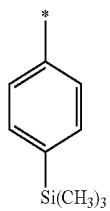


wherein in Formulae 1A-1 to 1A-3,
 ring A₁ is selected from a benzene group and a naphthalene group, ring A₂ is selected from a pyridine group, a quinoline group, and an isoquinoline group,
 X₁ is N(R₁₁)_{b11},
 R₁₁ is selected from groups represented by Formulae 6-1 to 6-27,
 b11 is 1,
 R₁ is selected from groups represented by Formulae 8-1 to 8-159,
 b1 is 1,
 R₃ to R₆, R₁₃, Z₁, and Z₂ 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 C₁-C₂₀ alkoxy group, a phenyl group, a biphenyl group, a terphenyl group, and a naphthyl group, and
 b3 to b6, b13, d4, d5, and d8 are each independently selected from 0, 1, and 2:



231
-continued

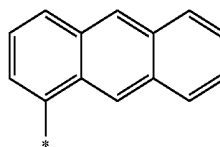


232

-continued

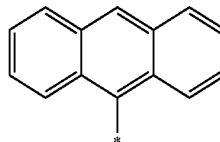
Formula 6-10

5

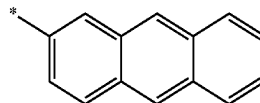


Formula 6-11

10



15



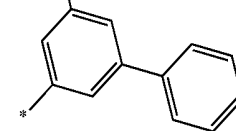
Formula 6-12

20



Formula 6-13

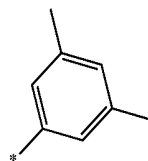
25



30

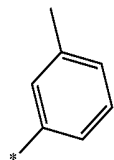
Formula 6-14

35



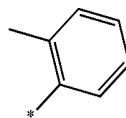
Formula 6-15

40



Formula 6-16

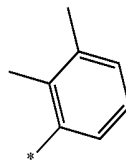
45



50

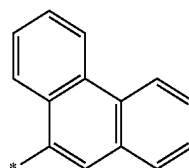
Formula 6-17

55

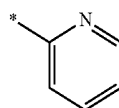


Formula 6-18

60



65



Formula 6-19

Formula 6-20

Formula 6-21

Formula 6-22

Formula 6-23

Formula 6-24

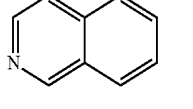
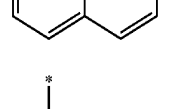
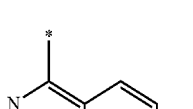
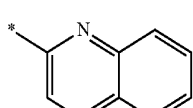
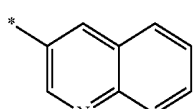
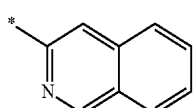
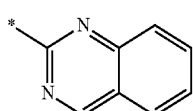
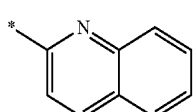
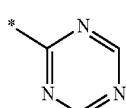
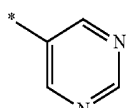
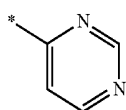
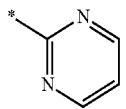
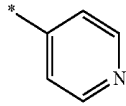
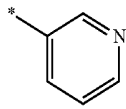
Formula 6-25

Formula 6-26

Formula 6-27

Formula 8-1

233
-continued



234
-continued

Formula 8-2

5

Formula 8-3

10

Formula 8-4

15

Formula 8-5

20

Formula 8-6

25

Formula 8-7

30

Formula 8-8

35

Formula 8-9

40

Formula 8-10

45

Formula 8-11

50

Formula 8-12

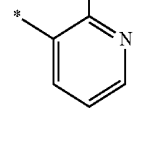
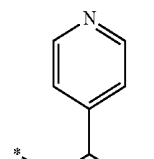
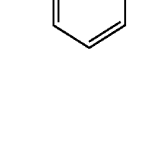
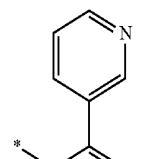
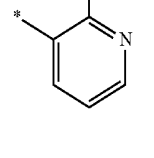
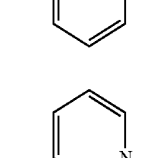
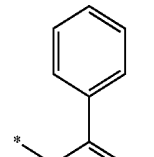
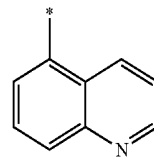
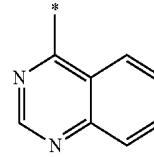
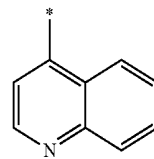
55

Formula 8-13

60

Formula 8-14

65



Formula 8-15

Formula 8-16

Formula 8-17

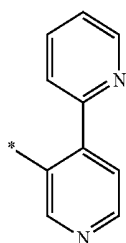
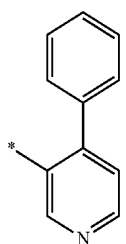
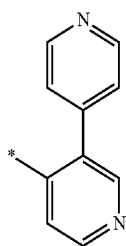
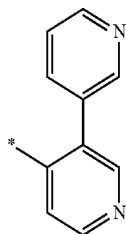
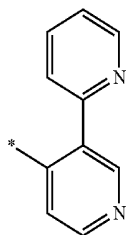
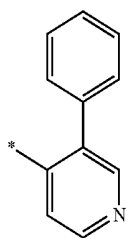
Formula 8-18

Formula 8-19

Formula 8-20

Formula 8-21

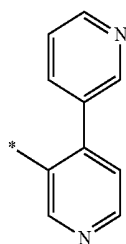
235
-continued



236
-continued

Formula 8-22

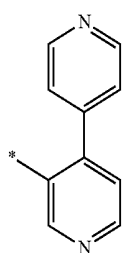
5



10

Formula 8-23

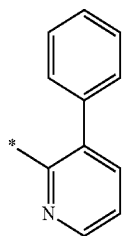
15



20

Formula 8-24

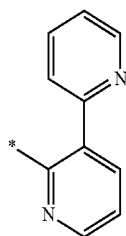
25



30

Formula 8-25

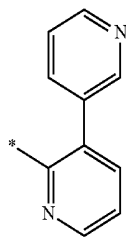
35



40

Formula 8-26

45

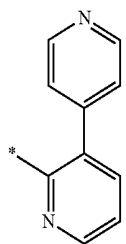


50

55

Formula 8-27

60



65

Formula 8-28

Formula 8-29

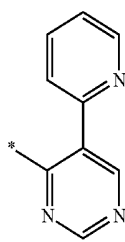
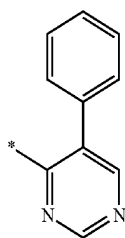
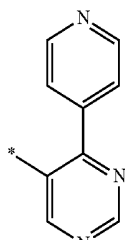
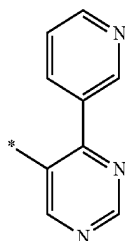
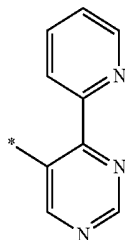
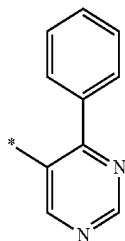
Formula 8-30

Formula 8-31

Formula 8-32

Formula 8-33

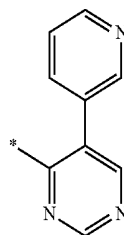
237
-continued



238
-continued

Formula 8-34

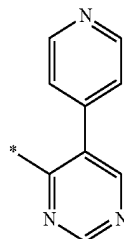
5



10

Formula 8-35

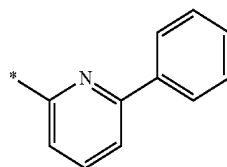
15



20

Formula 8-36

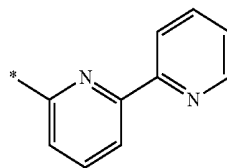
25



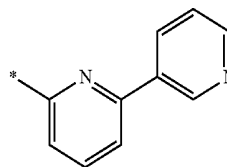
30

Formula 8-37

35

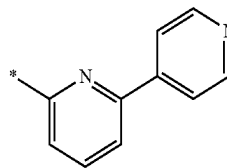


40

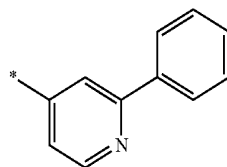


Formula 8-38

45



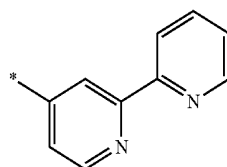
50



55

Formula 8-39

60



65

Formula 8-40

Formula 8-41

Formula 8-42

Formula 8-43

Formula 8-44

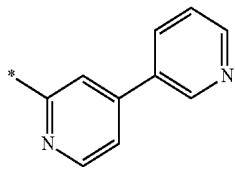
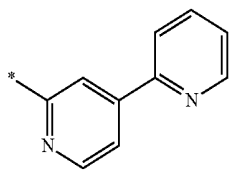
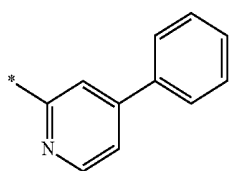
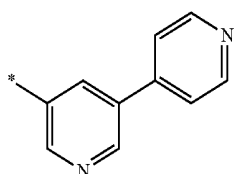
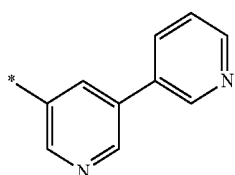
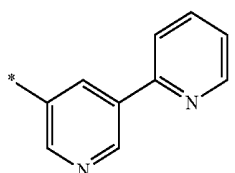
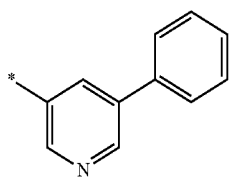
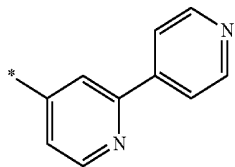
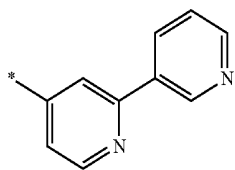
Formula 8-45

Formula 8-46

Formula 8-47

239

-continued

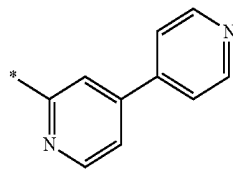


240

-continued

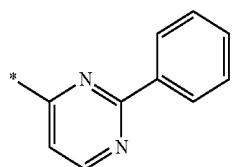
Formula 8-48

5



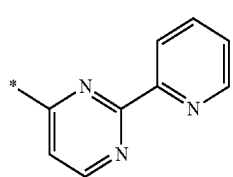
Formula 8-49

10



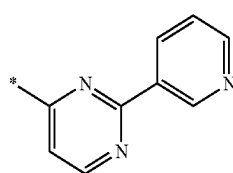
Formula 8-50

15



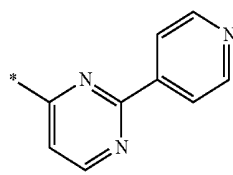
Formula 8-51

25



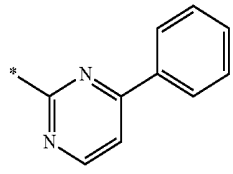
Formula 8-52

30



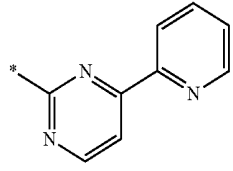
Formula 8-53

35



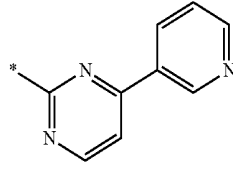
Formula 8-54

45



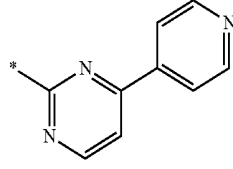
Formula 8-55

50



Formula 8-56

60



65

Formula 8-57

Formula 8-58

Formula 8-59

Formula 8-60

Formula 8-61

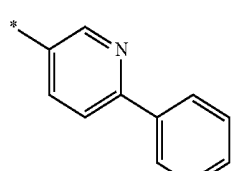
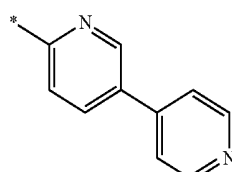
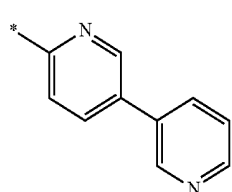
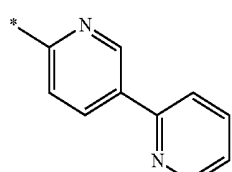
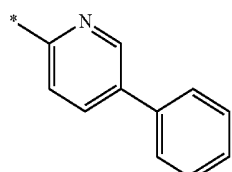
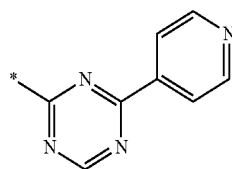
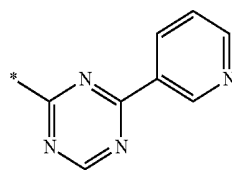
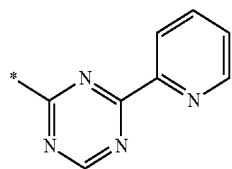
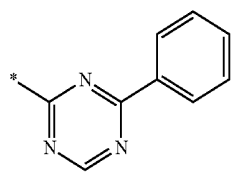
Formula 8-62

Formula 8-63

Formula 8-64

Formula 8-65

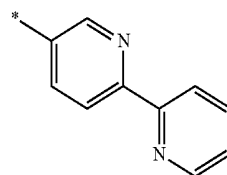
241
-continued



242
-continued

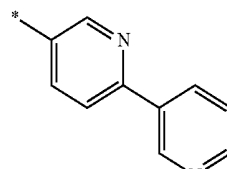
Formula 8-66

5



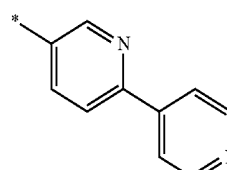
Formula 8-67

10



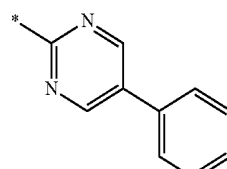
Formula 8-68

15



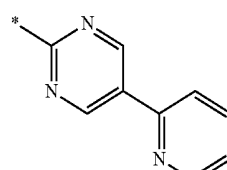
Formula 8-69

25



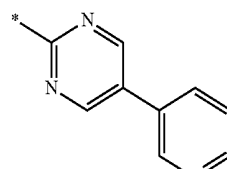
Formula 8-70

30



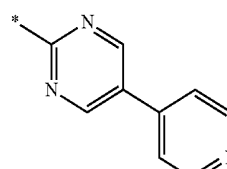
Formula 8-71

40



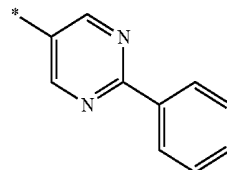
Formula 8-72

45



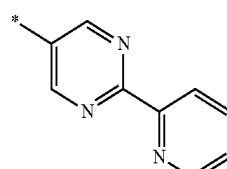
Formula 8-73

55



Formula 8-74

60



65

Formula 8-75

Formula 8-76

Formula 8-77

Formula 8-78

Formula 8-79

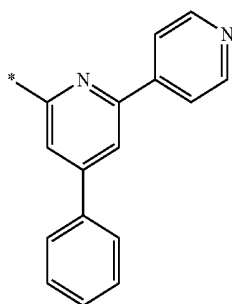
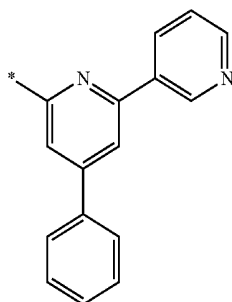
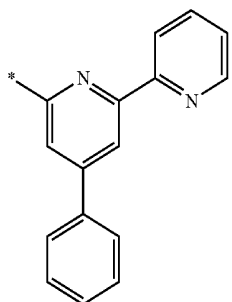
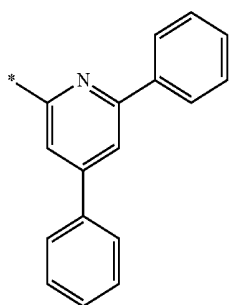
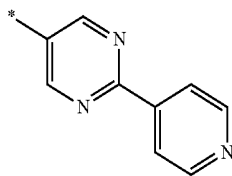
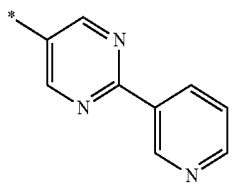
Formula 8-80

Formula 8-81

Formula 8-82

Formula 8-83

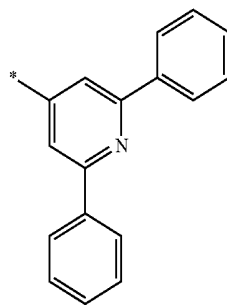
243
-continued



244
-continued

Formula 8-84

5



Formula 8-85 10

15

Formula 8-86

20

25

Formula 8-87

30

35

Formula 8-88

45

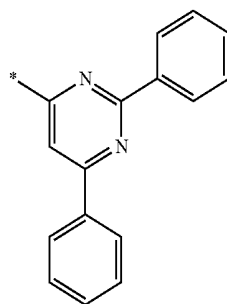
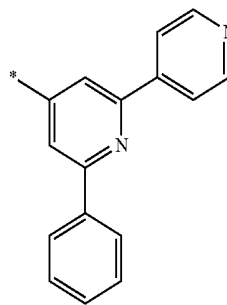
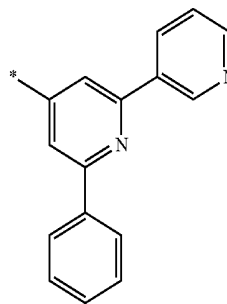
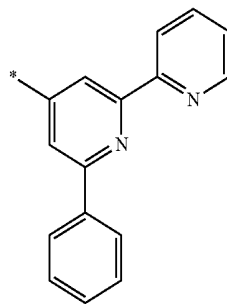
50

Formula 8-89

55

60

65



Formula 8-90

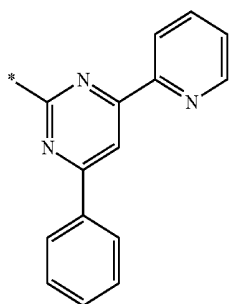
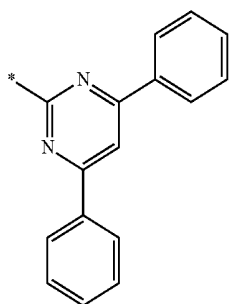
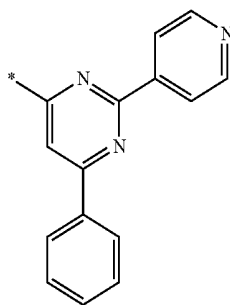
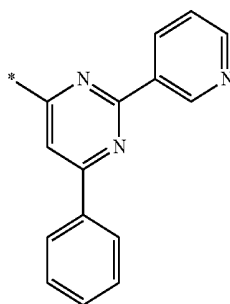
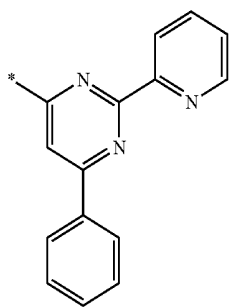
Formula 8-91

Formula 8-92

Formula 8-93

Formula 8-94

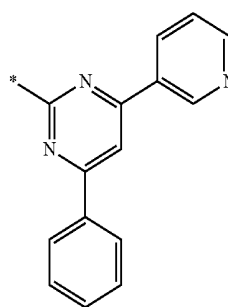
245
-continued



246
-continued

Formula 8-95

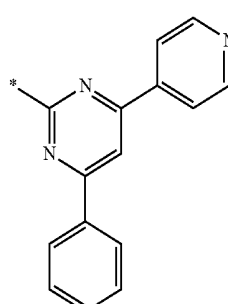
5



10

Formula 8-96

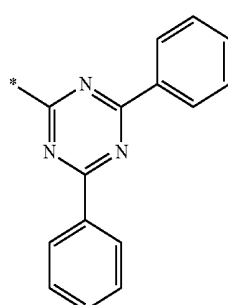
15



20

Formula 8-97

25

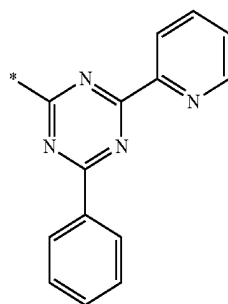


30

35

Formula 8-98

40

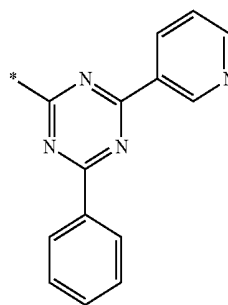


45

50

Formula 8-99

55



60

65

Formula 8-100

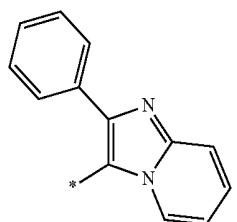
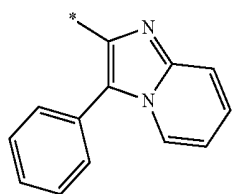
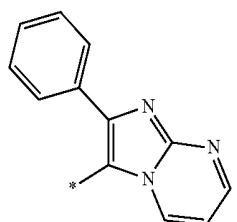
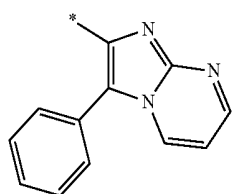
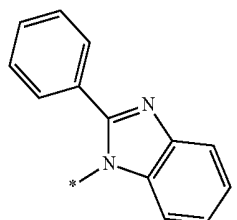
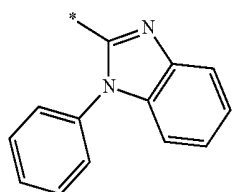
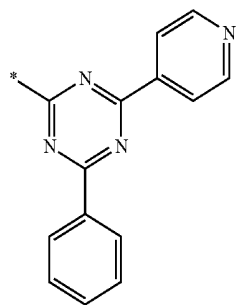
Formula 8-101

Formula 8-102

Formula 8-103

Formula 8-104

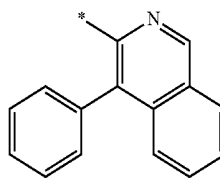
247
-continued



248
-continued

Formula 8-105

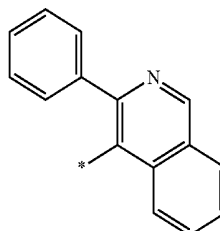
5



10

Formula 8-106

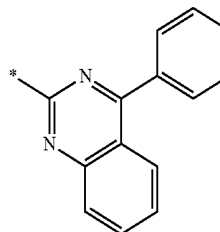
15



20

Formula 8-107

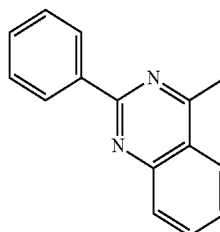
25



30

Formula 8-108

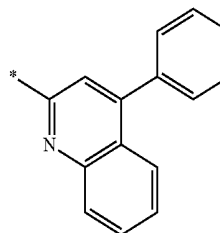
35



40

Formula 8-109

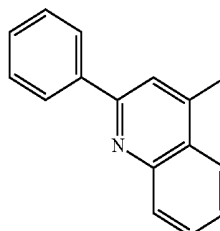
45



50

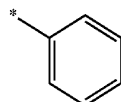
Formula 8-110

55

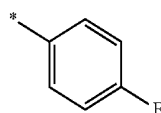


Formula 8-111

60



65



Formula 8-112

Formula 8-113

Formula 8-114

Formula 8-115

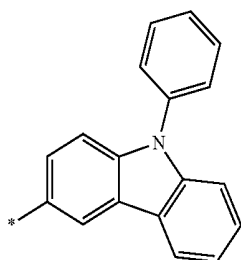
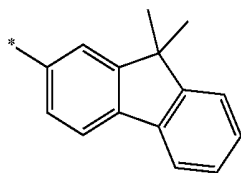
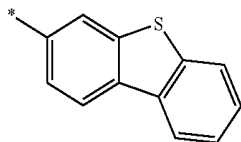
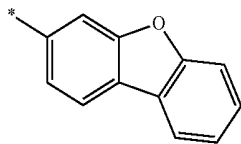
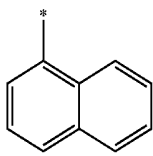
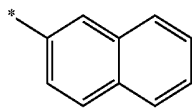
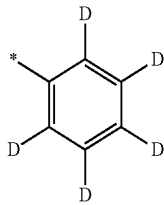
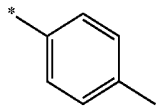
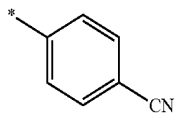
Formula 8-116

Formula 8-117

Formula 8-118

Formula 8-119

249
-continued



250
-continued

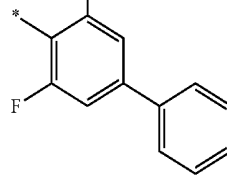
Formula 8-120

5



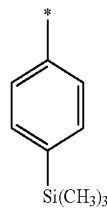
Formula 8-121

10



Formula 8-122

15



Formula 8-129

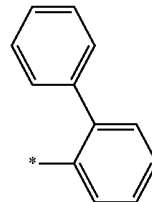
Formula 8-123

20

Formula 8-130

Formula 8-124

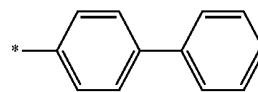
25



Formula 8-131

Formula 8-125

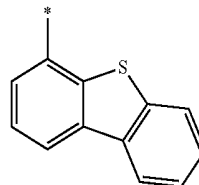
30



Formula 8-132

Formula 8-126

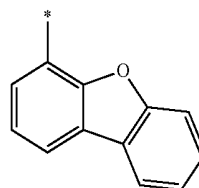
35



Formula 8-133

Formula 8-127

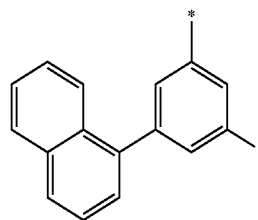
40



Formula 8-134

Formula 8-128

45



Formula 8-135

50

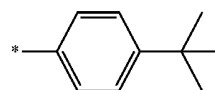
55

Formula 8-129

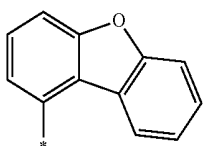
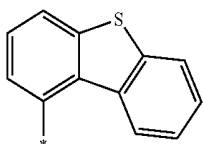
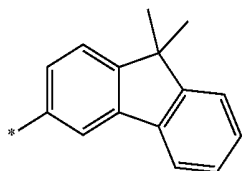
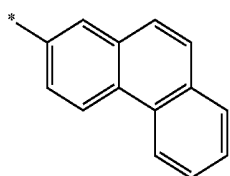
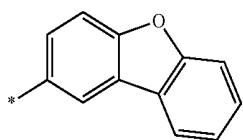
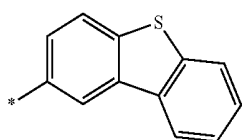
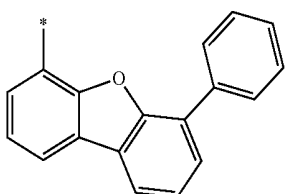
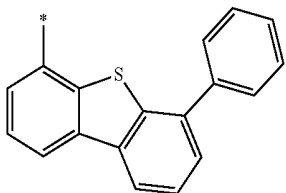
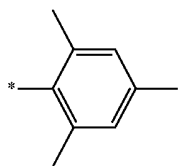
60

Formula 8-136

65



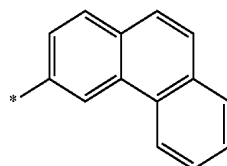
251
-continued



252
-continued

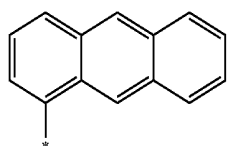
Formula 8-137

5



Formula 8-138

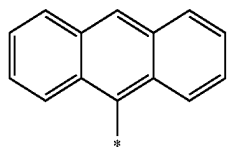
10



15

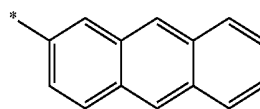
Formula 8-139

20



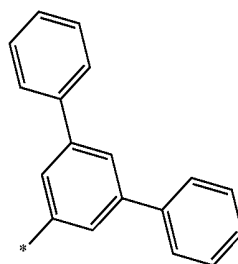
Formula 8-140

25



30

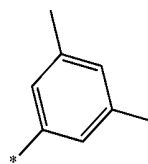
Formula 8-141



35

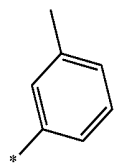
Formula 8-142

40



Formula 8-143

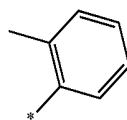
45



50

Formula 8-144

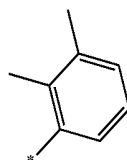
55



60

Formula 8-145

65



Formula 8-146

Formula 8-147

Formula 8-148

Formula 8-149

Formula 8-150

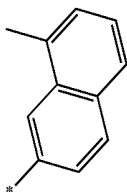
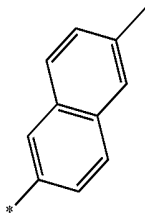
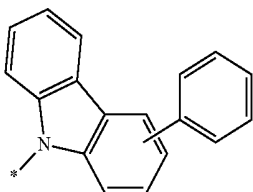
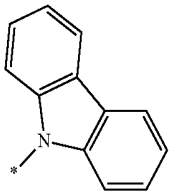
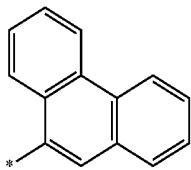
Formula 8-151

Formula 8-152

Formula 8-153

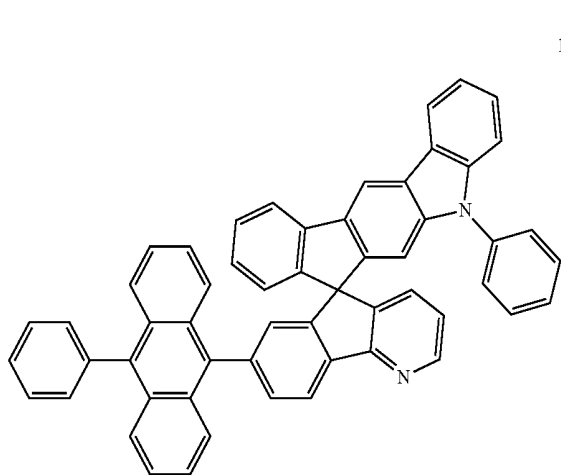
Formula 8-154

253
-continued



wherein, in Formulae 6-1 to 6-27 and 8-1 to 8-159, * indicates a binding site to an adjacent atom.

17. A condensed-cyclic compound represented by one of Compounds 1 to 35:



254
-continued

Formula 8-155

5

Formula 8-156

10

Formula 8-157

20

Formula 8-158

25

Formula 8-159

35

40

50

55

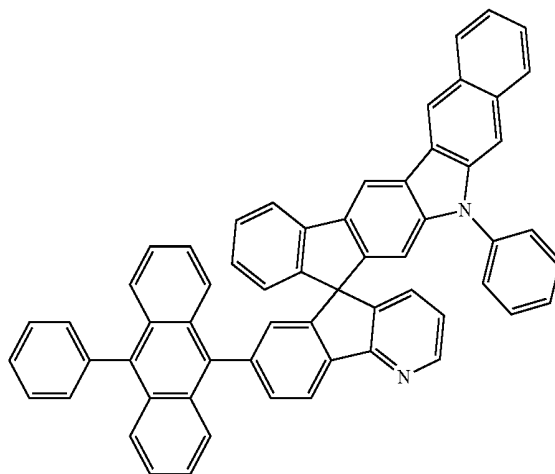
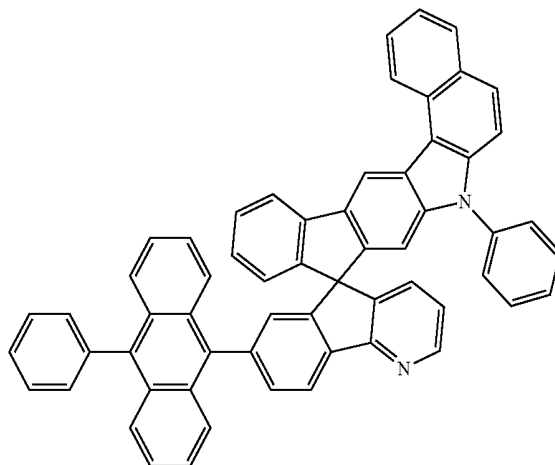
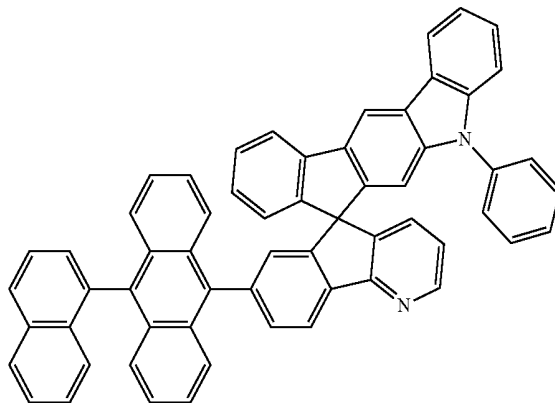
60

65

2

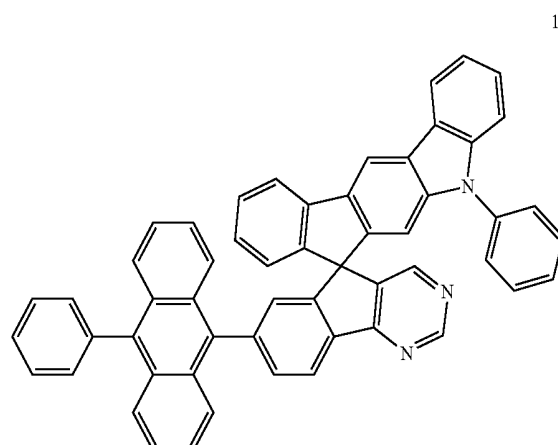
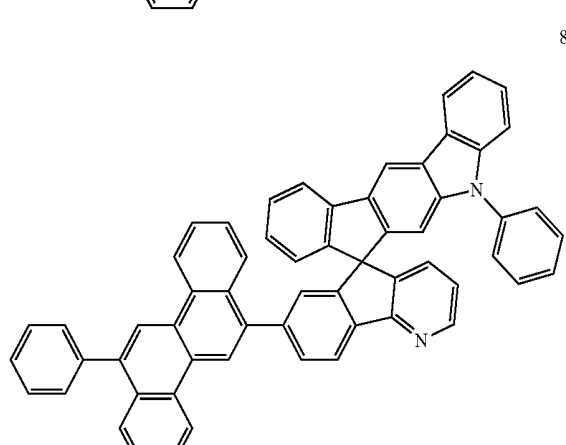
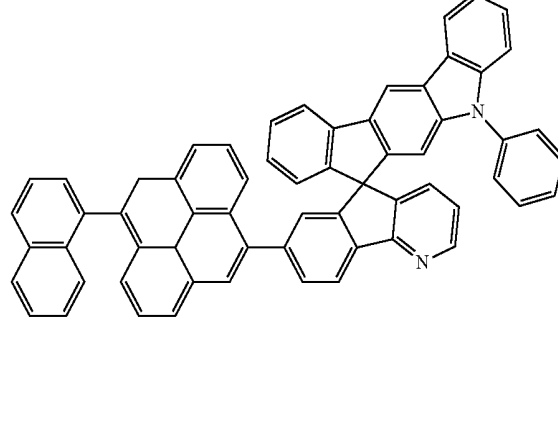
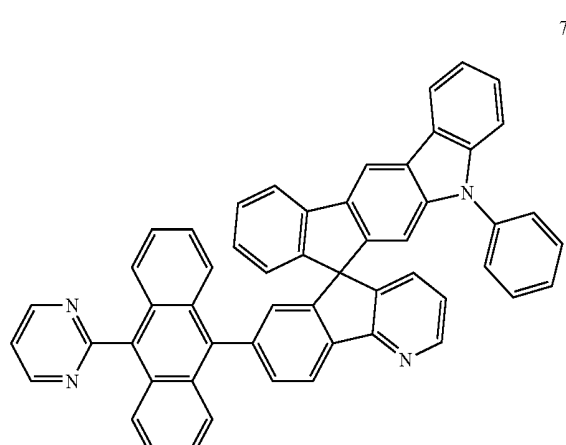
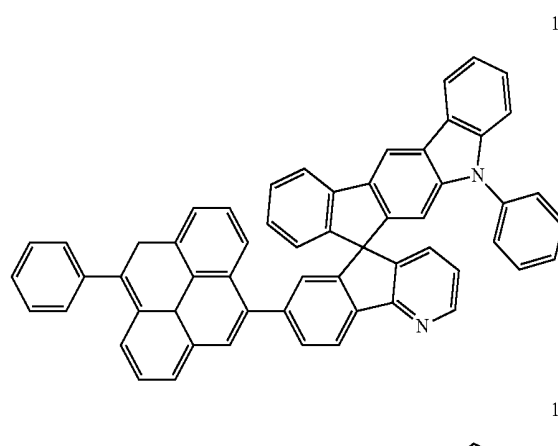
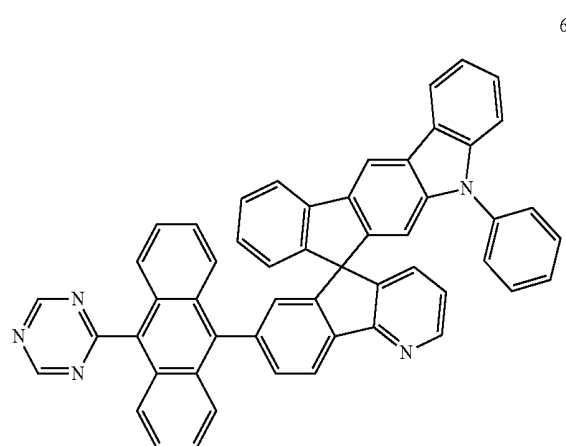
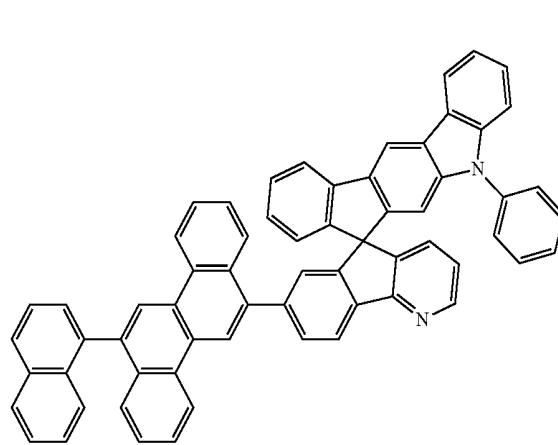
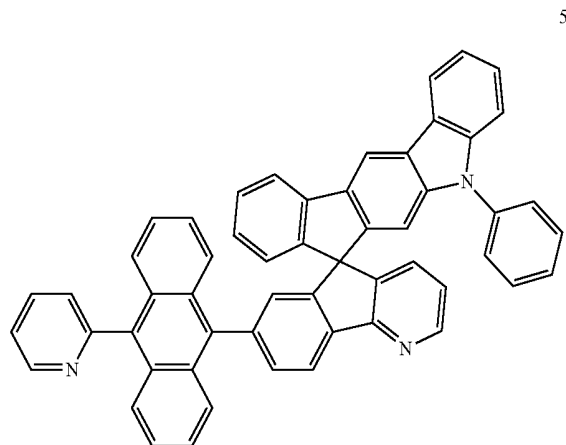
3

4



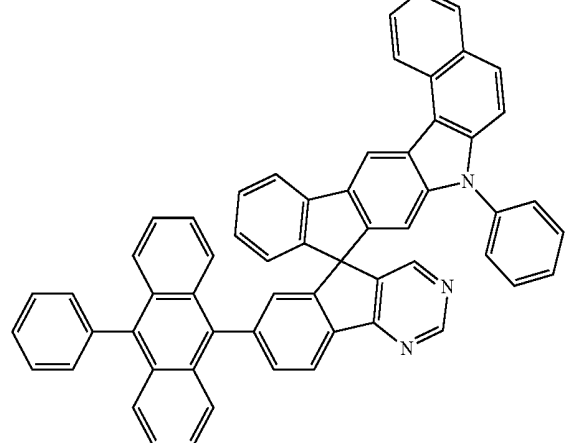
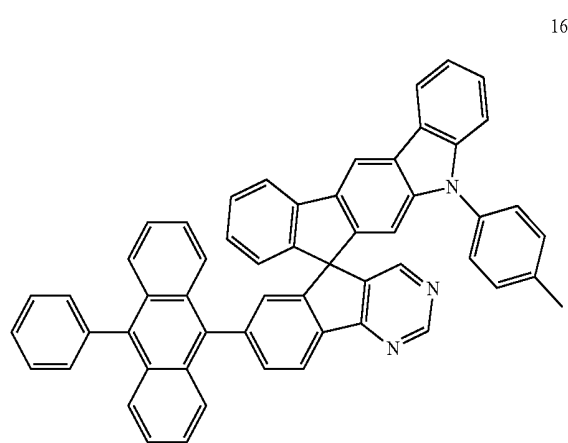
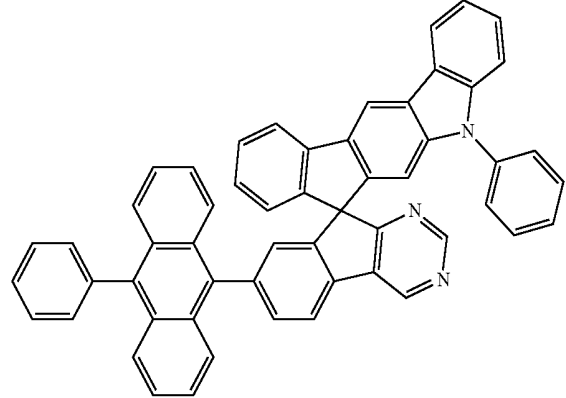
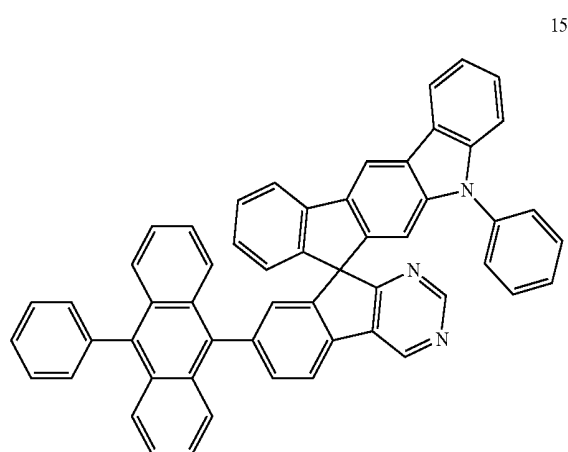
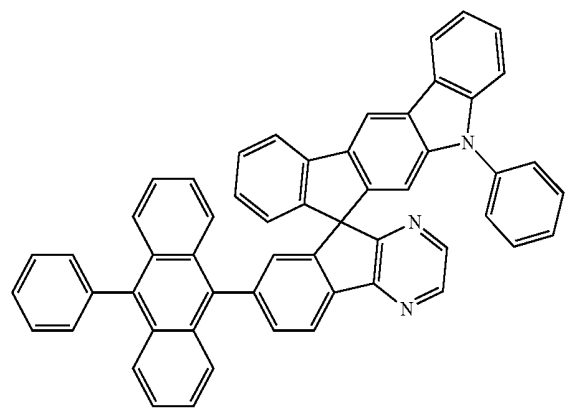
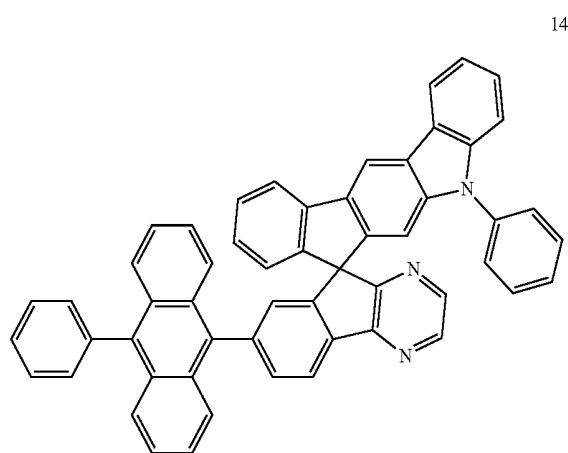
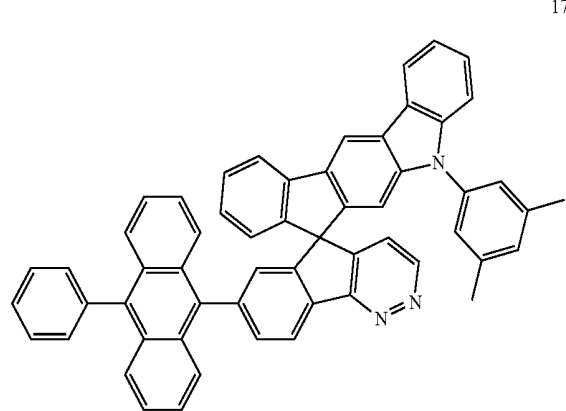
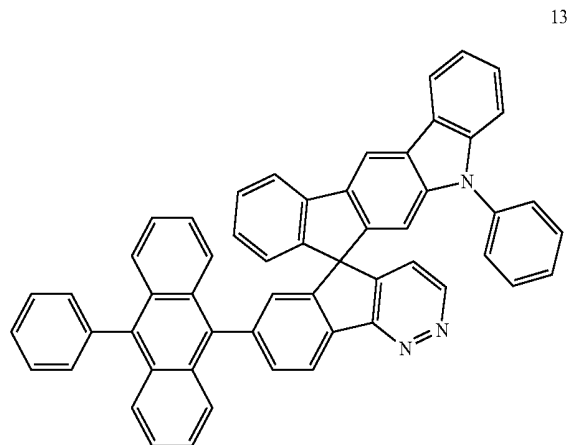
255
-continued

256
-continued

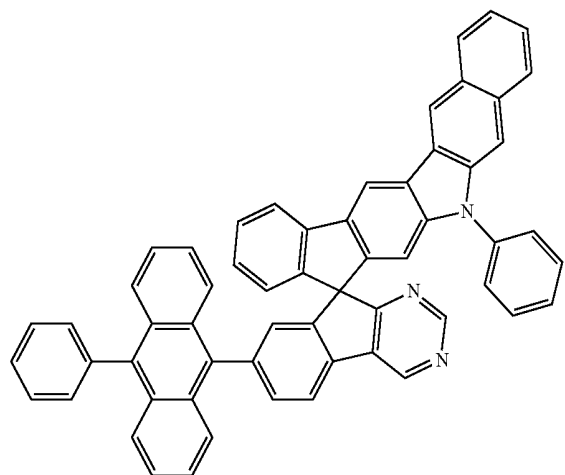
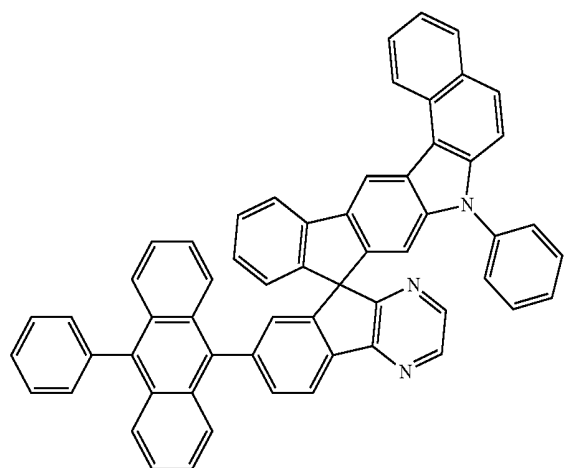
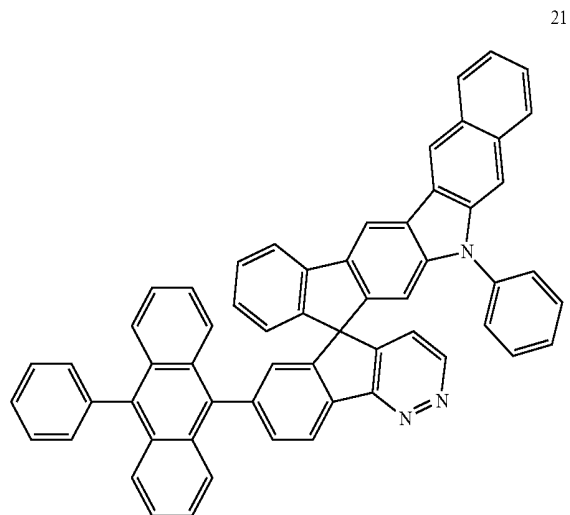


257
-continued

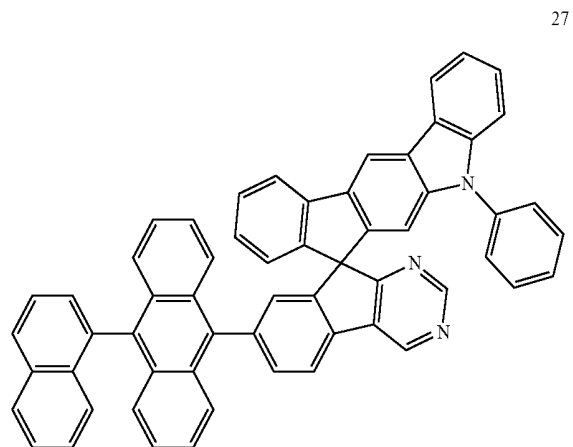
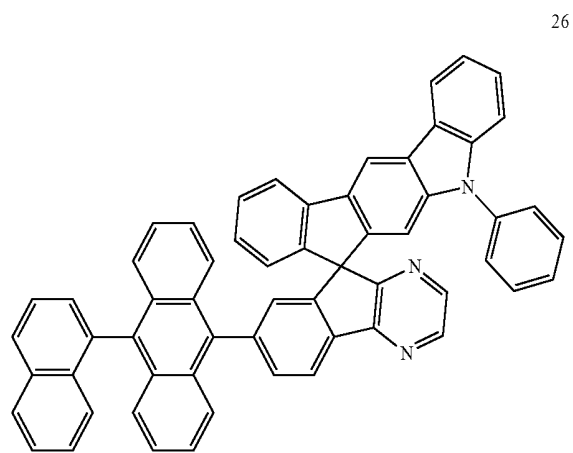
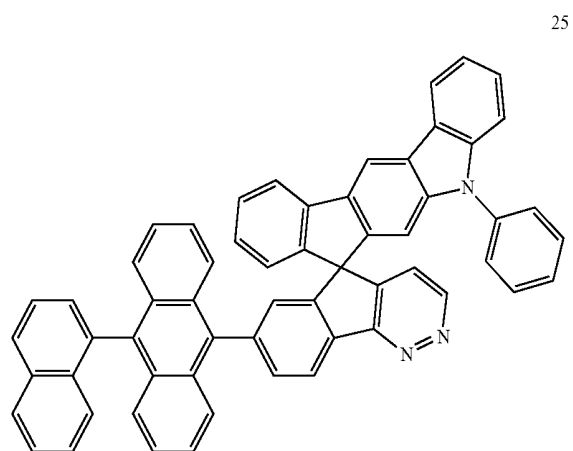
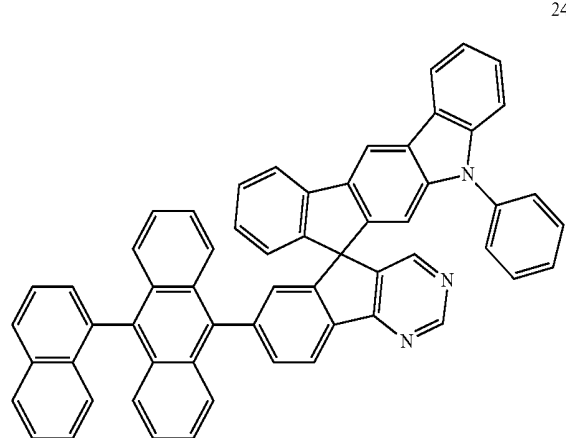
258
-continued



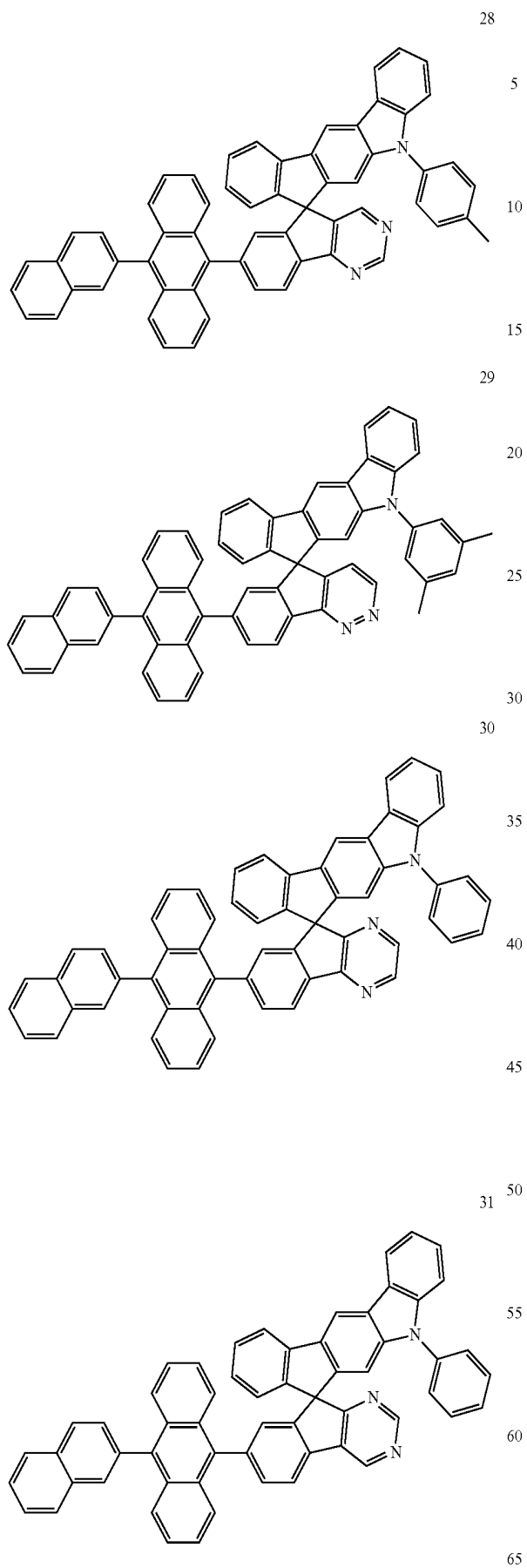
259
-continued



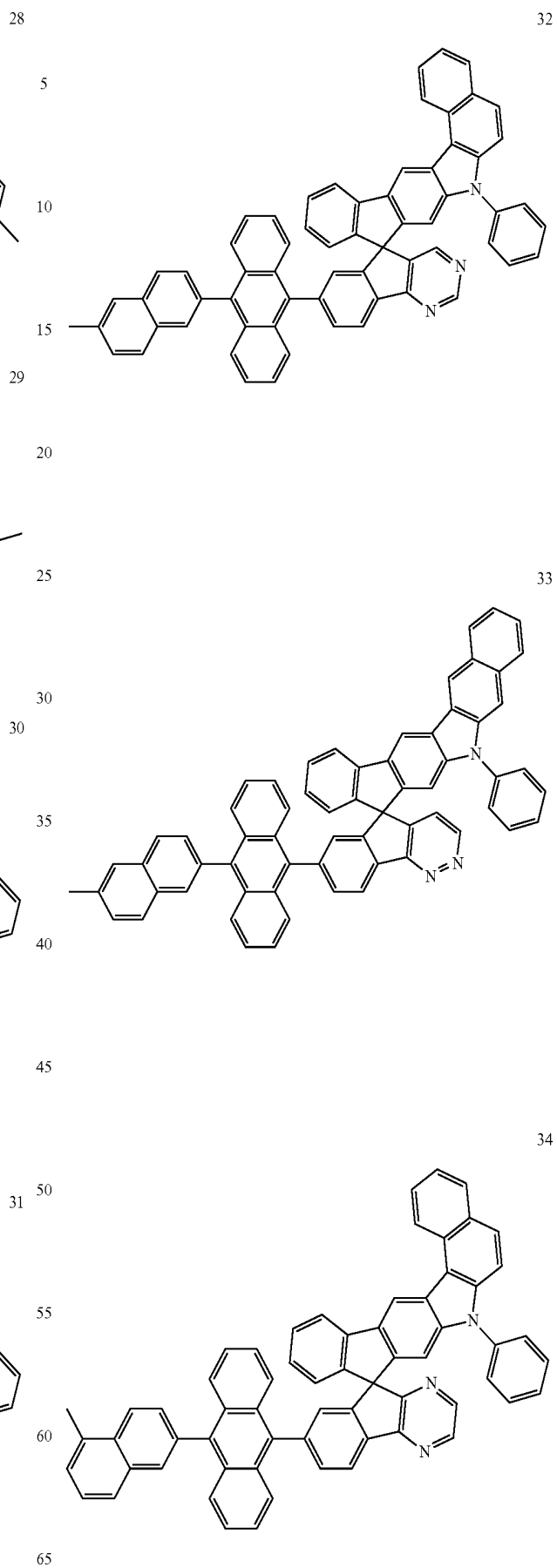
260
-continued



261
-continued

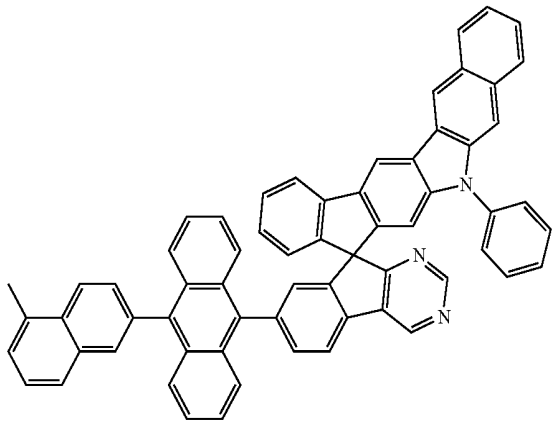


262
-continued



263

-continued



18. An organic light-emitting device comprising: a first electrode; a second electrode facing the first electrode; and an organic layer disposed between the first electrode and the

264

second electrode, wherein the organic layer comprises an emission layer and at least one condensed-cyclic compound of claim 1.

19. The organic light-emitting device of claim 18, wherein the first electrode is an anode,

the second electrode is a cathode,

the organic layer comprises i) a hole transport region between the first electrode and the emission layer that comprises at least one selected from a hole injection layer, a hole transport layer, an emission auxiliary layer, and an electron blocking layer and ii) an electron transport region between the emission layer and the

second electrode that comprises at least one selected from a hole blocking layer, an electron transport layer, and an electron injection layer, and

at least one selected from the hole transport region and the emission layer comprises the at least one condensed-cyclic compound.

20. The organic light-emitting device of claim 18, wherein the emission layer comprises the at least one condensed-cyclic compound.

* * * * *

专利名称(译)	稠环化合物和包括该稠环化合物的有机发光器件		
公开(公告)号	US10693076	公开(公告)日	2020-06-23
申请号	US15/703770	申请日	2017-09-13
[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	HWANG JINSOO KIM MYEONGSUK KIM SUNGWOOK		
发明人	HWANG, JINSOO KIM, MYEONGSUK KIM, SUNGWOOK		
IPC分类号	H01L51/00 C07F3/00 C07D307/77 H01L51/50 C07F9/6558 C07F9/655 C07D471/10 C07D307/91 C07D403/04 C07D405/10 C07D405/12 C07D405/14 C07D409/04 C07D409/14		
CPC分类号	C07F9/65517 H01L51/0073 C07D405/10 H01L51/5092 C07D409/14 H01L51/0052 H01L51/5056 H01L51/0072 C07D307/77 C07F9/65586 C07D405/12 C07D471/10 C07D409/04 C07D307/91 H01L51 /5096 C07D405/14 H01L51/5012 C07D403/04 H01L51/5072 C07F3/006 H01L51/0058 H01L51/0067 H01L2251/558 C07D487/10 H01L51/0054		
审查员(译)	LOEWE , ROBERT小号		
优先权	1020160118212 2016-09-13 KR		
其他公开文献	US20180076392A1		
外部链接	Espacenet		

摘要(译)

提供了稠环化合物和包括该稠环化合物的有机发光装置。稠环化合物由式1表示：其中，在式1、2A和2B中，环A 1选自苯基，萘基，吡啶基，噻吩基，吡嗪基，喹啉基，异喹啉基，喹喔啉基，喹唑啉基和cinnoline基。环A 2选自吡啶基，噻吩基，吡嗪基，喹啉基，异喹啉基，喹喔啉基，喹唑啉基和辛啉基；并响起A 3选自式2A表示的基团和式2B表示的基团。提供了有关化合物的更多细节。

10

190

150

110