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(54) Amine-based compound and organic light-emitting diode including the same

Aminbasierte Verbindung und organische lichtemittierende Diodenvorrichtung damit

Composé à base d'amine et diode électroluminescente organique le comprenant

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US-A- 5 935 721

- **KUM HEE LEE ET AL: "Synthesis and Electroluminescent Properties of Blue Fluorescent Triphenylamine Substituted Anthracene Derivatives for OLEDs", MOLECULAR CRYSTALS AND LIQUID CRYSTALS, vol. 530, no. 1, 11 October 2010 (2010-10-11), pages 48/[204]-55/[211], XP55060418, ISSN: 1542-1406, DOI: 10.1080/15421406.2010.495882**

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Description

[0001] The present invention relates to an amine-based compound for organic light-emitting diodes, and an organic light-emitting diode including the compound.

[0002] Organic light-emitting diodes (OLEDs), which are self-emitting diodes, have advantages such as wide viewing angles, excellent contrast, quick response, high brightness, excellent driving voltage characteristics, and can provide multicoloured images.

[0003] A typical OLED has a structure including a substrate, and an anode, a hole transport layer (HTL), an emission layer (EML), an electron transport layer (ETL), and a cathode which are sequentially stacked on the substrate. In this regard, the HTL, the EML, and the ETL are organic thin-films comprising organic compounds.

[0004] An operating principle of an OLED having the above-described structure is as follows.

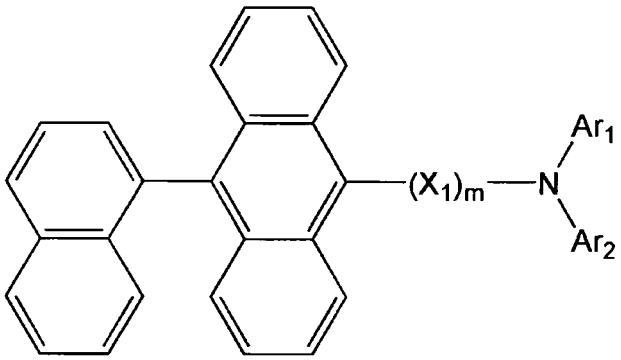
[0005] When a voltage is applied between the anode and the cathode, holes injected from the anode move to the EML via the HTL, and electrons injected from the cathode move to the EML via the ETL. The holes and electrons recombine in the EML to generate excitons. When the excitons drop from an excited state to a ground state, light is emitted.

[0006] Kum Hee Lee et.al. disclose blue fluorescent triphenylamine substituted anthracenes derivatives, especially 9-(naphthalen-2-yl)-10-(4-triphenylamine)anthracene, comprised in the organic layer of organic light-emitting devices ("Synthesis and electroluminescent properties of blue fluorescent triphenylamine substituted anthracene derivatives for OLEDs", Molecular Crystals and Liquid Crystals, Vol. 530, No.1, 11 October 2010)

[0007] The present invention provides an amine-based compound having a novel structure and an organic light-emitting diode including the amine-based compound.

[0008] According to a first aspect of the present invention there is provided an amine-based compound represented by Formula 1 below:

Formula 1



wherein, in Formula 1, Ar_1 and Ar_2 are each independently a substituted or unsubstituted C_6-C_{60} aryl group or a substituted or unsubstituted C_2-C_{60} heteroaryl group;

X_1 is a substituted or unsubstituted C_6-C_{60} arylene group or a substituted or unsubstituted C_2-C_{60} heteroarylene group;

m is an integer from 1 to 5; and

at least one substituent of each of the substituted C_6-C_{60} aryl group, the substituted C_2-C_{60} heteroaryl group, the substituted C_6-C_{60}arylene group, and the substituted C_2-C_{60} heteroarylene group is one of a deuterium atom; -F; -Cl; -Br; -I; -CN; a hydroxyl group; $-NO_2$; an amino group; an amidino group; hydrazine; hydrazone; a carboxyl group or a salt thereof; a sulfonic acid group or a salt thereof; a phosphoric acid or a salt thereof; a tri(C_6-C_{60} aryl)silyl group; a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group; a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group and a C_2-C_{60} alkynyl group that is substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof; a C_3-C_{60} cycloalkyl group, a C_3-C_{60} cycloalkenyl group, a C_6-C_{60} aryl group, a C_2-C_{60} heteroaryl group, a C_6-C_{60} aralkyl group, a C_6-C_{60} aryloacy group, a C_6-C_{60} arylthiol group; a C_3-C_{60} cycloalkyl group, a C_3-C_{60} cycloalkenyl group, a C_6-C_{60} aryl group, a C_2-C_{60} heteroaryl group, a C_6-C_{60} aralkyl group, a C_6-C_{60} aryloxy group, and a C_6-C_{60} arylthiol group that is substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic

acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one fluorine (F), a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group and a C_2-C_{60} heteroaryl group,

wherein at least one of Ar_1 and Ar_2 is a C_6-C_{60} aryl group substituted with at least one electron withdrawing group selected from the group consisting of -F; -CN; -NO_2; a C_1-C_{60} alkyl group substituted with at least one -F; a C_2-C_{60} heteroaryl group; and a C_2-C_{60} heteroaryl group substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one -F, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group, and a C_2-C_{60} heteroaryl group.

[0009] According to another aspect of the present invention, there is provided an organic light-emitting diode comprising a first electrode, a second electrode disposed opposite to the first electrode, and an organic layer disposed between the first electrode and the second electrode, the organic layer comprising at least one of the amine-based compounds according to the invention in its first aspect.

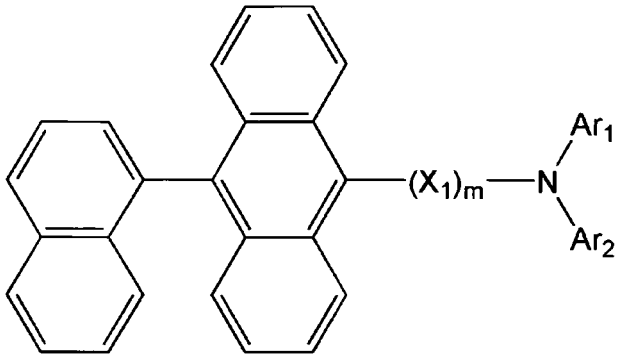
[0010] A more complete appreciation of the present invention, and many of the attendant advantages thereof, will be readily apparent as the same becomes better understood by reference to the following detailed description when considered in conjunction with the accompanying drawings, wherein:

Figure. 1 schematically illustrates the structure of an organic light-emitting diode according to an embodiment of the invention.

[0011] As used herein, the term "and/or" includes any and all combinations of one or more of the associated listed items.

[0012] According to an aspect of the present invention, there is provided an amine-based compound represented by Formula 1:

Formula 1



[0013] In Formula 1 above, Ar_1 and Ar_2 are each independently a substituted or unsubstituted C_6-C_{60} aryl group, or a substituted or unsubstituted C_2-C_{60} heteroaryl group; and X_1 is a substituted or unsubstituted C_6-C_{60} arylene group, or a substituted or unsubstituted C_2-C_{60} heteroarylene group and m is an integer selected from 1, 2, 3, 4 and 5.

[0014] At least one substituent of each of the substituted C_6-C_{60} aryl group, the substituted C_2-C_{60} heteroaryl group, the substituted C_6-C_{60}arylene group, and the substituted C_2-C_{60} heteroarylene group is one of a deuterium atom; -F; -Cl; -Br; -I; -CN; a hydroxyl group; -NO_2; an amino group; an amidino group; hydrazine; hydrazone; a carboxyl group or a salt thereof; a sulfonic acid group or a salt thereof; a phosphoric acid or a salt thereof; a tri(C_6-C_{60}aryl)silyl group; a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, and a C_2-C_{60} alkynyl group; a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group and a C_2-C_{60} alkynyl group that is substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof; a C_3-C_{60} cycloalkyl group, a C_3-C_{60} cycloalkenyl group, a C_6-C_{60} aryl group, a C_2-C_{60} heteroaryl group, a C_6-C_{60} aralkyl group, a C_6-C_{60} aryloxy group, a C_6-C_{60} arylthiol group; a C_3-C_{60} cycloalkyl group, a C_3-C_{60} cycloalkenyl group, a C_6-C_{60} aryl group, a C_2-C_{60} heteroaryl group, a C_6-C_{60} aralkyl group, a C_6-C_{60} aryloxy group, and a C_6-C_{60} arylthiol group that is substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one fluorine (F), a

C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group and a C_2-C_{60} heteroaryl group.

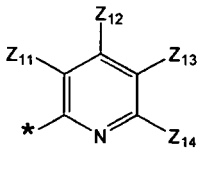
[0015] In Formula 1 above, at least one of Ar_1 and Ar_2 is a C_6-C_{60} aryl group substituted with at least one electron withdrawing group selected from the group consisting of -F; -CN; - NO_2; a C_1-C_{60} alkyl group substituted with at least one -F; a C_2-C_{60} heteroaryl group; and a C_2-C_{60} heteroaryl group substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one -F, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group, and a C_2-C_{60} heteroaryl group.

[0016] For example, the at least one electron withdrawing group may be selected from the group consisting of: -F; -CN; -NO_2; a C_1-C_{20} alkyl group substituted with at least one -F; a C_2-C_{20} heteroaryl group including a ring which contains a N atom; and a C_2-C_{20} heteroaryl group that includes a ring which contains a N atom and is substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a C_6-C_{20} aryl group, and a C_2-C_{20} heteroaryl group.

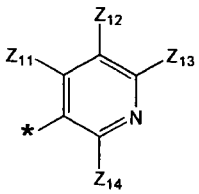
[0017] In an embodiment of the present invention, the at least one electron withdrawing group in Formula 1 is selected from the group consisting of -F; -CN; a C_1-C_{20} alkyl group substituted with at least one -F; a pyrrolyl group, a pyrazolyl group, an imidazolyl group, an imidazolinyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a benzoimidazolyl group, an indolyl group, a purinyl group, a quinolinyl group, an isoquinolinyl group, a phthalazinyl group, an indolizinyl group, a quinazolinyl group, a cinnolinyl group, an indazolyl group, a carbazolyl group, a phenazinyl group, a phenanthridinyl group, a triazinyl group, a pyridazinyl group, a triazoly group, and a tetrazoly; and a pyrrolyl group, a pyrazolyl group, an imidazolyl group, an imidazolinyl group, an imidazopyridinyl group, imidazopyrimidinyl, pyridinyl, pyrazinyl, pyrimidinyl, benzoimidazolyl, indolyl, purinyl, quinolinyl, isoquinolinyl , phthalazinyl, indolizinyl, quinazolinyl, cinnolinyl, indazolyl, carbazolyl, phenazinyl, phenanthridinyl, triazinyl, pyridazinyl, triazolyl, and a tetrazolyl group that is substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group and a carbazolyl group.

[0018] For example, the at least one electron withdrawing group in Formula 1 above may be selected from, but is not limited to, the group consisting of: -F; -CN; a C_1-C_{20} alkyl group substituted with at least one -F; a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group; a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group that is substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group.

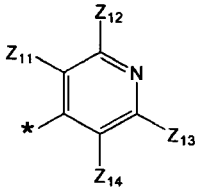
[0019] In an embodiment of the invention, the at least one electron withdrawing group in Formula 1 above is selected from the group consisting of -F; -CN; -CH_2F; -CHF_2; -CF_3; and groups represented by Formulae 2(1) to 2(14) below:



Formula 2(1)



Formula 2(2)



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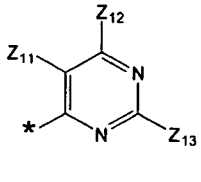

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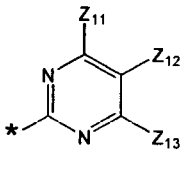

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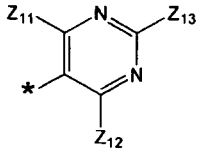

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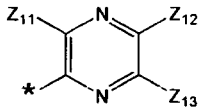

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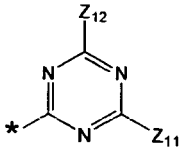

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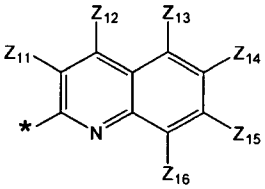

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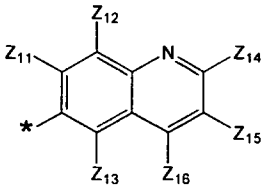

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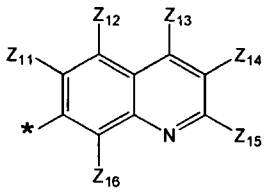

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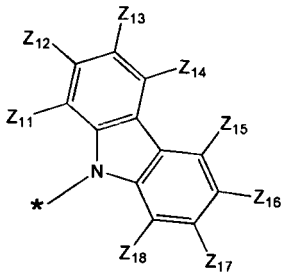

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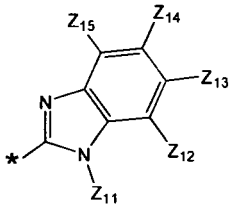
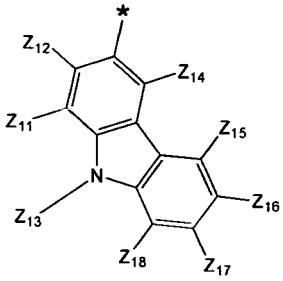


Formula 2(13)



Formula 2(14)

[0020] In Formulae 2(1) to 2(14) above, Z_{11} to Z_{18} may each independently be a hydrogen atom, a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, or a carbazolyl group.

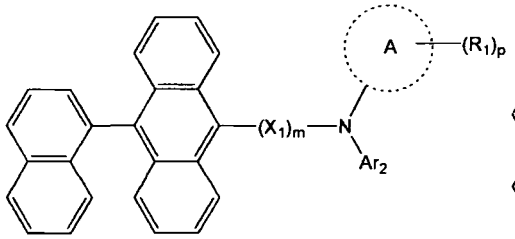
[0021] For example, Z_{11} to Z_{18} in Formulae 2(1) to 2(14) above may each independently be a hydrogen atom, a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonate group or a salt thereof, a phosphoric acid or a salt thereof, a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, a methoxy group, an ethoxy group, a propoxy group, a butoxy group, a pentoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, or a carbazolyl group, but are not limited thereto.

[0022] In Formula 1 above, the at least one of Ar_1 and Ar_2 may be a C_6-C_{60} aryl group substituted with at least two electron withdrawing groups.

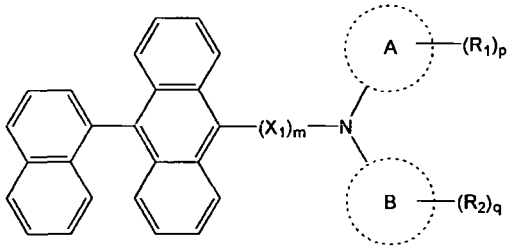
[0023] In an embodiment of the present invention, the at least one of Ar_1 and Ar_2 is a phenyl group, a biphenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, or a fluorenyl group that is substituted with at least two electron withdrawing groups. The electron withdrawing groups may each independently be selected from the group consisting of a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; and a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a phthalazinyl group, a benzoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN; a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group.

[0024] At least one of Ar_1 and Ar_2 in the amine-based compound of Formula 1 above is a C_6-C_{60} aryl group substituted with at least one of the above electron withdrawing groups. Accordingly, the amine-based compound may be represented by Formula 1(1) or Formula 1(2) below:

Formula 1(1)



Formula 1(2)



[0025] In Formula 1(1) above, Ar_2 is a substituted or unsubstituted C_6-C_{20} aryl group, or a substituted or unsubstituted C_2-C_{20} heteroaryl group. In Formulae 1(1) and 1(2) above, the rings A and B are each independently a substituted C_6-C_{20} aryl group; R_1 and R_2 are each independently an electron withdrawing group selected from the group consisting of: -F; -CN; -NO_2; a C_1-C_{60} alkyl group substituted with at least one -F; a C_2-C_{60} heteroaryl group; and a C_2-C_{60} heteroaryl group substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one -F, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group, and a C_2-C_{60} heteroaryl group; and p and q are each independently an integer selected from 1, 2, 3, 4, 5, 6, 7, 8 and 9.

[0026] The electron withdrawing group in Formulae 1(1) and 1(2) are as described above, and thus a detailed description thereof will not be repeated herein.

[0027] For example, the amine-based compound may be represented by Formula 1(1) above, wherein at least one of R_1 in Formula 1(1) may be -CN.

[0028] In an embodiment of the present invention, the amine-based compound is represented by Formula 1(2) above, wherein at least one of R_1 and R_2 in Formula 1(2) is -CN.

[0029] In an embodiment of the present invention, the amine-based compound is represented by Formula 1(1) above, wherein the ring A in Formula 1(1) is a substituted phenyl group, a substituted biphenyl group, a substituted naphthyl group, a substituted anthryl group, a substituted phenanthrenyl group, a substituted pyrenyl group, or a substituted fluorenyl group, but is not limited thereto.

[0030] In another embodiment of the present invention, the amine-based compound is represented by Formula 1(2) above, wherein the rings A and B in Formula 1(2) may be each independently a substituted phenyl group, a substituted biphenyl group, a substituted naphthyl group, a substituted anthryl group, a substituted phenanthrenyl group, a substituted pyrenyl group, or a substituted fluorenyl group.

[0031] The amine-based compound may be represented by Formula 1(1), wherein the A ring may be a substituted phenyl group, a substituted biphenyl group, a substituted naphthyl group, a substituted anthryl group, a substituted phenanthrenyl group, a substituted pyrenyl group, or a substituted fluorenyl group; R_1 may be at least one electron withdrawing group selected from the group consisting of a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; and a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a phthalazinyl group, a benzoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN; a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group; and p may be 2, 3, or 4, for example, may be 2.

[0032] In an embodiment of the present invention, the amine-based compound is represented by Formula 1(2), wherein the ring A and the ring B are each independently a substituted phenyl group, a substituted biphenyl group, a substituted naphthyl group, a substituted anthryl group, a substituted phenanthrenyl group, a substituted pyrenyl group, or a substituted fluorenyl group; R_1 and R_2 are each independently at least one electron withdrawing group selected from the group consisting of a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; and a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a phthalazinyl group, a benzoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN; a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group; and p and q are each

independently an integer of 2, 3, or 4, for example, p and q may both be an integer of 2.

[0033] At least one of Ar_1 and Ar_2 in Formula 1 above may be a phenyl group, a biphenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, or a fluorenyl group that are substituted with at least one of the above-listed electron withdrawing groups.

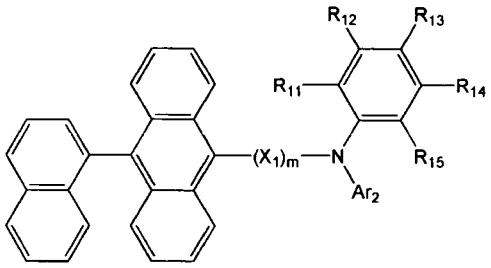
[0034] In Formula 1 above, Ar_1 and Ar_2 may each independently be a substituted or unsubstituted phenyl group, a substituted or unsubstituted pentalenyl group, a substituted or unsubstituted indenyl group, a substituted or unsubstituted naphthyl group, a substituted or unsubstituted azulenyl group, a substituted or unsubstituted heptalenyl group, a substituted or unsubstituted indacenyl group, a substituted or unsubstituted acenaphthyl group, a substituted or unsubstituted fluorenyl group, a substituted or unsubstituted phenalenyl group, a substituted or unsubstituted phenanthrenyl group, a substituted or unsubstituted anthryl group, a substituted or unsubstituted fluoranthenyl group, a substituted or unsubstituted triphenylenyl group, a substituted or unsubstituted pyrenyl group, a substituted or unsubstituted chrysenyl group, a substituted or unsubstituted naphthacenyl group, a substituted or unsubstituted picenyl group, a substituted or unsubstituted perylenyl group, a substituted or unsubstituted pentaphenyl group, a substituted or unsubstituted hexacenyl group, a substituted or unsubstituted pyrrolyl group, a substituted or unsubstituted pyrazolyl group, a substituted or unsubstituted imidazolyl group, a substituted or unsubstituted imidazolinyl group, a substituted or unsubstituted imidazopyridinyl group, a substituted or unsubstituted imidazopyrimidinyl group, a substituted or unsubstituted pyridinyl group, a substituted or unsubstituted pyrazinyl group, a substituted or unsubstituted pyrimidinyl group, a substituted or unsubstituted benzoimidazolyl group, a substituted or unsubstituted indolyl group, a substituted or unsubstituted purinyl group, a substituted or unsubstituted quinolinyl group, a substituted or unsubstituted phthalazinyl group, a substituted or unsubstituted indolizinyl group, a substituted or unsubstituted naphthyridinyl group, a substituted or unsubstituted quinazolinyl group, a substituted or unsubstituted cinnolinyl group, a substituted or unsubstituted indazolyl group, a substituted or unsubstituted carbazolyl group, a substituted or unsubstituted phenazinylene group, a substituted or unsubstituted phenanthridinyl group, a substituted or unsubstituted pyranyl group, a substituted or unsubstituted chromenyl group, a substituted or unsubstituted furanyl group, a substituted or unsubstituted benzofuranyl group, a substituted or unsubstituted thiophenyl group, a substituted or unsubstituted benzothiophenyl group, a substituted or unsubstituted isothiazolyl group, a substituted or unsubstituted benzoimidazolyl group, a substituted or unsubstituted isoxazolyl group, a substituted or unsubstituted dibenzothiophenyl group, a substituted or unsubstituted dibenzofuranyl group, a substituted or unsubstituted triazinyl group, a substituted or unsubstituted oxadiazolyl group, a substituted or unsubstituted pyridazinyl group, a substituted or unsubstituted triazolyl group, a substituted or unsubstituted tetrazolyl group, or a substituted or unsubstituted phenanthrolinyl group. At least one of Ar_1 and Ar_2 may be a phenyl group, a biphenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, or a fluorenyl group that are substituted with at least one of the above-listed electron withdrawing groups.

[0035] For example, Ar_1 and Ar_2 in Formula 1 above may each independently be a substituted or unsubstituted phenyl group, a substituted or unsubstituted naphthyl group, a substituted or unsubstituted fluorenyl group, a substituted or unsubstituted phenanthrenyl group, a substituted or unsubstituted anthryl group, a substituted or unsubstituted triphenylenyl group, a substituted or unsubstituted pyrenyl group, a substituted or unsubstituted chrysenyl group, a substituted or unsubstituted pyridinyl group, a substituted or unsubstituted pyrazinyl group, a substituted or unsubstituted pyrimidinyl group, a substituted or unsubstituted quinolinyl group, a substituted or unsubstituted carbazolyl group, a substituted or unsubstituted triazinyl group, a substituted or unsubstituted dibenzothiophenyl group, a substituted or unsubstituted dibenzofuranyl group, or a substituted or unsubstituted phenanthrolinyl group. At least one of Ar_1 and Ar_2 may be a phenyl group, a biphenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, or a fluorenyl group that is substituted with at least one electron withdrawing group selected from the group consisting of -F; -CN; a C_1-C_{20} alkyl group substituted with at least one -F; a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; and a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group.

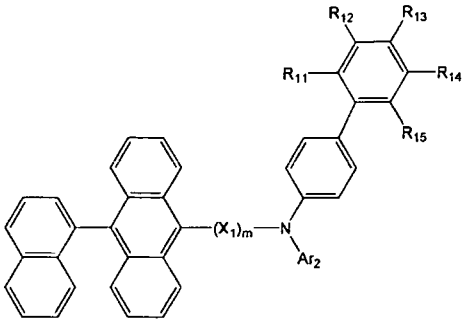
[0036] In Formula 1, Ar_1 and Ar_2 may be linked together by a single bond.

[0037] The amine-based compound may be represented by any one of Formulae 1A to iJ below:

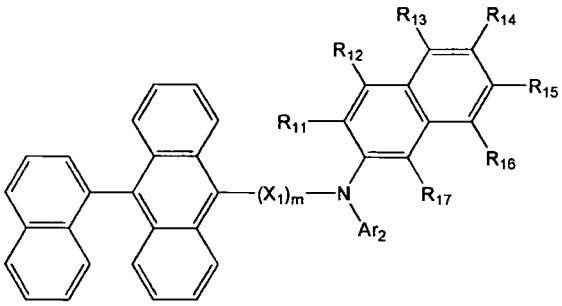
Formula 1A



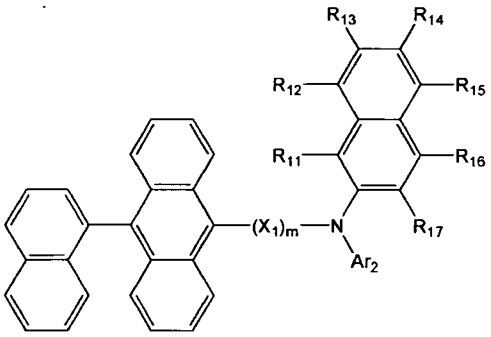
Formula 1B



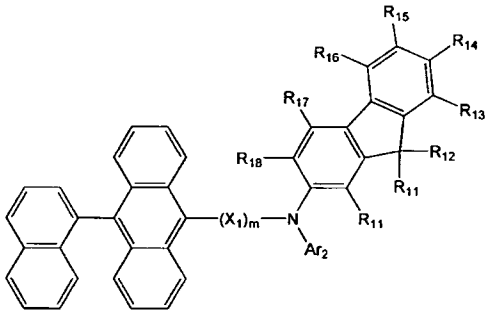
Formula 1C



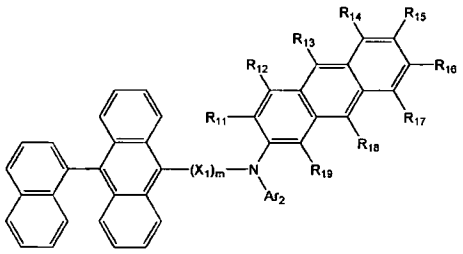
Formula 1D



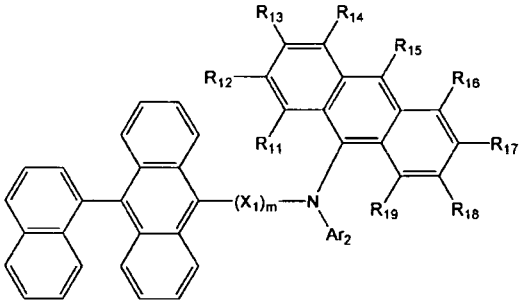
Formula 1E



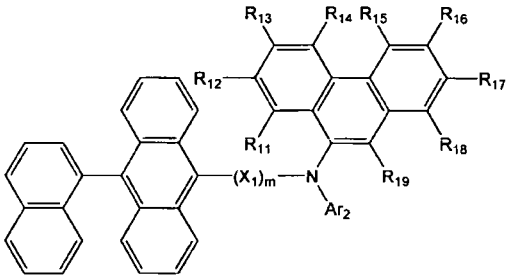
Formula 1F



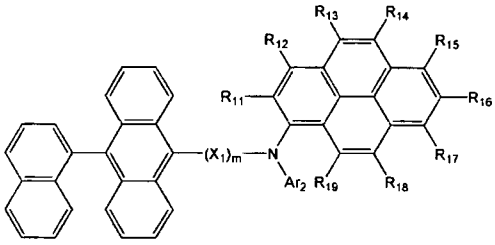
Formula 1G



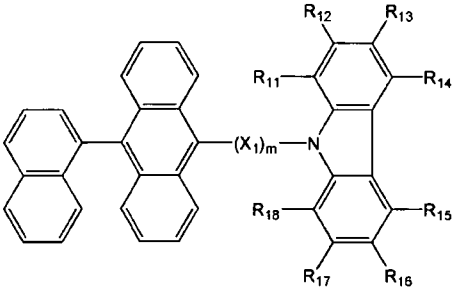
Formula 1H



Formula 1I



Formula 1J



Ar_2 in Formulae 1A to 1I may be the same as described above in conjunction with other formulae.

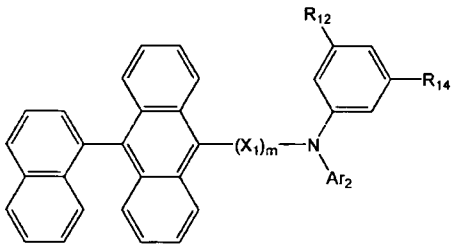
[0038] At least one of R_{11} to R_{15} in Formulae 1A and 1B, at least one of R_{11} to R_{17} in Formulae 1C and 1D, at least one of R_{11} to R_{18} in Formulae 1E and 1J, and at least one of R_{11} to R_{19} in Formula 1F, 1G, 1H and 1I may each independently be an electron withdrawing group selected from the group consisting of -F; -CN; $-NO_2$; a C_1-C_{60} alkyl group substituted with at least one -F; a C_2-C_{60} heteroaryl group; and a C_2-C_{60} heteroaryl group substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one -F, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group, and a C_2-C_{60} heteroaryl group.

[0039] For example, at least one of R_{11} to R_{15} of Formulae 1A and 1B, at least one of R_{11} to R_{17} in Formulae 1C and 1D, at least one of R_{11} to R_{18} of Formulae 1E and 1J, and at least one of R_{11} to R_{19} of Formulae 1F, 1G, 1H and 1I may each independently be, but are not limited to, an electron withdrawing group selected from the group consisting of: -F; -CN; a C_1-C_{20} alkyl group substituted with at least one -F; a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a phthalazinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, - Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group.

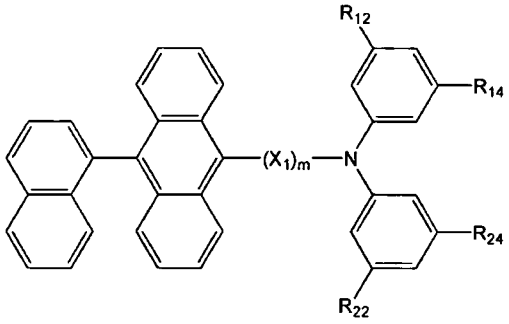
[0040] In an embodiment of the present invention, Ar_2 in Formulae 1A to 1I may be, but are not limited thereto, a phenyl group, a biphenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, or a fluorenyl group that are substituted with at least one electron withdrawing group selected from the group consisting of -F; -CN; a C_1-C_{20} alkyl group substituted with at least one -F; a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; and a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a phthalazinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group.

[0041] In an embodiment of the invention, the amine-based compound may be represented by Formula 1A-(1) or 1A-(2) below:

Formula 1A-(1)



Formula 1A-(2)



[0042] In Formulae 1A-(1) and 1A-(2), R_{12}, R_{14}, R_{22}, and R_{24} may each independently be an electron withdrawing group selected from the group consisting of a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; and a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a phthalazinyl group, a benzoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN; a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group; and Ar_2 may be a substituted or unsubstituted phenyl group, a substituted or unsubstituted biphenyl group, a substituted or unsubstituted naphthyl group, a substituted or unsubstituted anthryl group, a substituted or unsubstituted phenanthrenyl group, a substituted or unsubstituted pyrenyl group, or a substituted or unsubstituted fluorenyl group.

[0043] In Formulae 1A-(1) and 1A-(2), R_{12}, R_{14}, R_{22} and R_{24} may each independently be selected from the group consisting of -F; -CN; $-CH_2F$; $-CHF_2$; $-CF_3$; and groups represented by Formulae 2(1) to 2(14) above.

[0044] In Formulae 1A-(1) and 1A-(2) above, R_{12}, R_{14}, R_{22} and R_{24} may each independently be one of the groups represented by Formulae 2(1) to 2(8) above.

[0045] In Formulae 1A-(1) and 1A-(2) above, R_{12}, R_{14}, R_{22} and R_{24} may each independently be the group represented by Formula 2(2) above, but are not limited thereto.

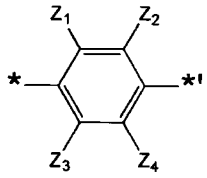
[0046] In Formula 1A-(1), Ar_2 may be a phenyl group, a biphenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, or a fluorenyl group, but are not limited thereto.

[0047] In Formula 1, X_1 may be a substituted or unsubstituted phenylene group, a substituted or unsubstituted pentalenylene group, a substituted or unsubstituted indenylene group, a substituted or unsubstituted naphthylene group, a substituted or unsubstituted azulenylene group, a substituted or unsubstituted heptalenylene group, a substituted or unsubstituted indacenylene group, a substituted or unsubstituted acenaphthylene group, a substituted or unsubstituted fluorenylene group, a substituted or unsubstituted phenalenylene group, a substituted or unsubstituted phenanthrenylene group, a substituted or unsubstituted anthrylene group, a substituted or unsubstituted fluoranthenylene group, a substituted or unsubstituted triphenylenylene group, a substituted or unsubstituted pyrenylene group, a substituted or unsubstituted chrysenylene group, a substituted or unsubstituted naphthacenylene group, a substituted or unsubstituted picenylene group, a substituted or unsubstituted perylenylene group, a substituted or unsubstituted pentaphenylene group, a substituted or unsubstituted hexacenylene group, a substituted or unsubstituted pyrrolylene group, a substituted or unsubstituted pyrazolylene group, a substituted or unsubstituted imidazolylene group, a substituted or unsubstituted imidazolinylene group, a substituted or unsubstituted imidazopyridinylene group, a substituted or unsubstituted imidazopyrimidinylene group, a substituted or unsubstituted pyridinylene group, a substituted or unsubstituted pyrazinylene group, a substituted or unsubstituted pyrimidinylene group, a substituted or unsubstituted indolylene group, a substituted or unsubstituted purinylene group, a substituted or unsubstituted quinolinylene group, a substituted or unsubstituted phthalazinylene group, a substituted or unsubstituted indolizinylene group, a substituted or unsubstituted naphthyridinylene group, a substituted or unsubstituted quinazolinylene group, a substituted or unsubstituted cinnolinylene group, a substituted or unsubstituted indazolylene group, a substituted or unsubstituted carbazolylene group, a substituted or unsubstituted phenazinylene group, a substituted or unsubstituted phenanthridinylene group, a substituted or unsubstituted pyranylene group, a substituted or unsubstituted chromenylene group, a substituted or unsubstituted furanylene group, a substituted or unsubstituted benzofuranylene group, a substituted or unsubstituted thiophenylene group, a substituted or unsubstituted benzothiophenylene group, a substituted or unsubstituted isothiazolylene group, a substituted or unsubstituted benzoimidazolylene group, a substituted or unsubstituted isoxazolylene group, a substituted or unsubstituted dibenzothiophenylene group, a substituted or unsubstituted dibenzofuranylene group, a substituted or unsubstituted triazinylene group, a substituted or unsubstituted oxadiazolylene group, a substituted or unsubstituted

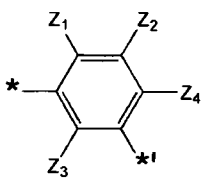
pyridazinylene group, a substituted or unsubstituted triazolylene group, or a substituted or unsubstituted tetrazolylene group. X_1 in Formula 1 may have at least one substituent, which may be selected from among the substituents described above.

[0048] In an embodiment, X_1 in Formula 1 is, but is not limited to, a substituted or unsubstituted phenylene group, a substituted or unsubstituted naphthylene group, a substituted or unsubstituted fluorenylene group, a substituted or unsubstituted phenanthrenylene group, a substituted or unsubstituted anthrylene group, a substituted or unsubstituted triphenylenylene group, a substituted or unsubstituted pyrenylene group, a substituted or unsubstituted chrysenylene group, a substituted or unsubstituted pyridinylene group, a substituted or unsubstituted pyrazinylene group, a substituted or unsubstituted pyrimidinylene group, a substituted or unsubstituted quinolinylene group, a substituted or unsubstituted quinazolinylene group, a substituted or unsubstituted carbazolylene group, a substituted or unsubstituted dibenzothiophenylene group, a substituted or unsubstituted dibenzofuranylene group, a substituted or unsubstituted triazinylene group, a substituted or unsubstituted pyridazinylene group, a substituted or unsubstituted triazolylene group, or a substituted or unsubstituted tetrazolylene group.

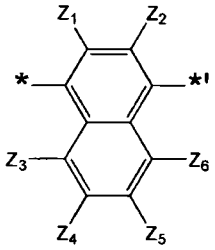
[0049] For example, X_1 in Formula 1 above may be a group represented by one of Formulae 5(1) to 5(16) below:



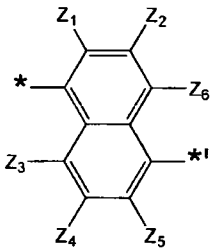
Formula 5(1)



Formula 5(2)



Formula 5(3)



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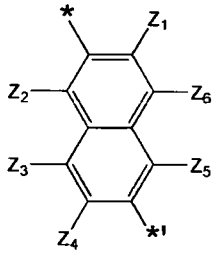

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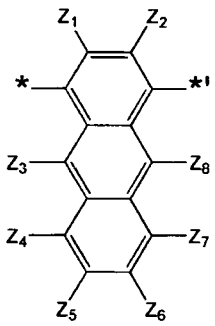

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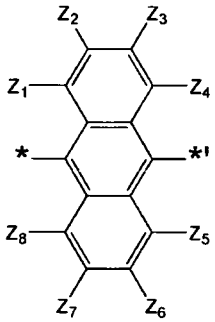

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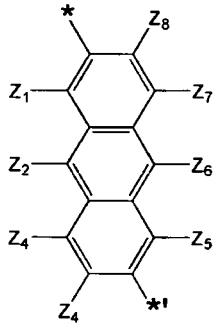

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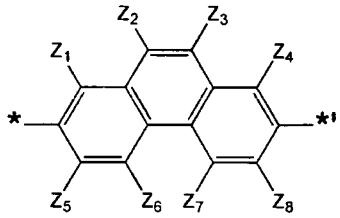

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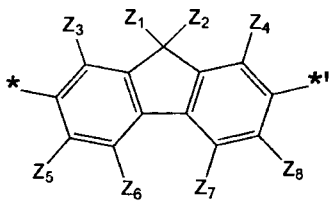

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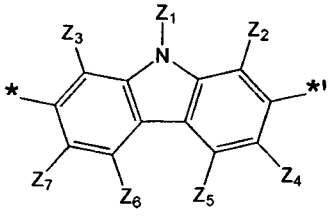

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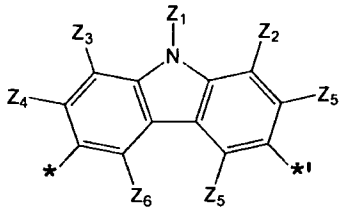

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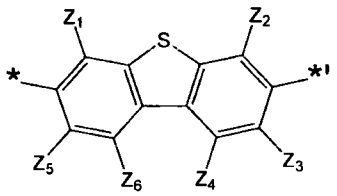

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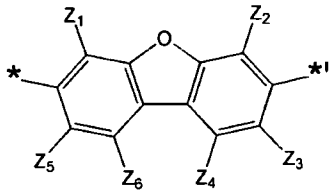

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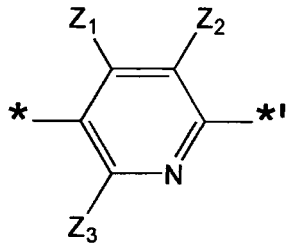

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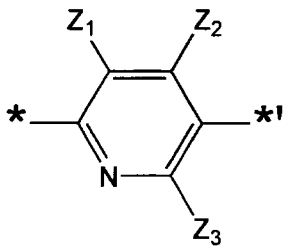


Formula 5(16)

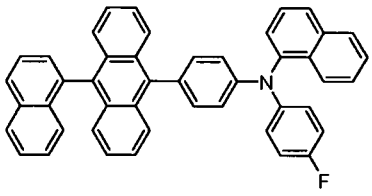
[0050] In Formulae 5(1) to 5(16), Z_1 to Z_8 may each independently be one of a hydrogen atom; a deuterium atom; -F; -Cl; -Br; -I; -CN; a hydroxyl group; $-NO_2$; an amino group; an amidino group; hydrazine; hydrazone; a carboxyl group or a salt thereof; a sulfonic acid group or a salt thereof; phosphoric acid or a salt thereof; a C_1-C_{20} alkyl group; a C_1-C_{20} alkoxy group; a C_1-C_{20} alkyl group and a C_1-C_{20} alkoxy groups that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, sulfonic acid group or a salt thereof, and phosphoric acid or a salt thereof; a C_6-C_{20} aryl group; a C_2-C_{20} heteroaryl group; and a C_6-C_{20} aryl group and a C_2-C_{20} heteroaryl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkoxy group, a C_6-C_{20} aryl group and a C_2-C_{20} heteroaryl group.

[0051] In Formulae 5(1) to 5(16), * indicates a binding site to anthracene in Formula 1, and *, indicates a binding site to N in Formula 1.

[0052] For example, Z_1 to Z_8 in Formulae 5(1) to 5(16) may each independently be, but are not limited to, a hydrogen atom; a deuterium atom; -F; -Cl; -Br; -I; -CN; a hydroxyl group; - NO_2; an amino group; an amidino group; hydrazine; hydrazone; a carboxyl group or a salt thereof; a sulfonic acid group or a salt thereof; a phosphoric acid or a salt thereof; a methyl group, an ethyl group, a propyl group, a butyl group, and a pentyl group; a methoxy group, an ethoxy group, a propoxy group, a butoxy group, and a pentoxy group; a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, a methoxy group, an ethoxy group, a propoxy group, a butoxy group, and a pentoxy group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, and a phosphoric acid or a salt thereof; a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, and a fluorenyl group; a pyridinyl group, a pyrimidinyl group, a triazinyl group, a quinolyl group, and a carbazolyl group; a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a fluorenyl group, a pyridinyl group, a pyrimidinyl group, a triazinyl group, a quinolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group and a pyridinyl group.

[0053] In Formula 1, m is an integer from 1 to 5. In an embodiment of the present invention, m is 1, 2, or 3, but is not limited thereto.

[0054] The amine-based compound of Formula 1 may be, for example, any one of Compounds 1 to 109 below, but is not limited thereto:



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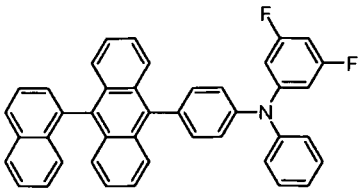

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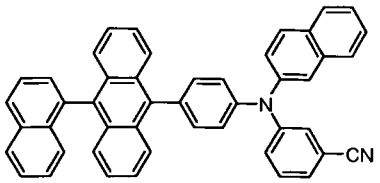

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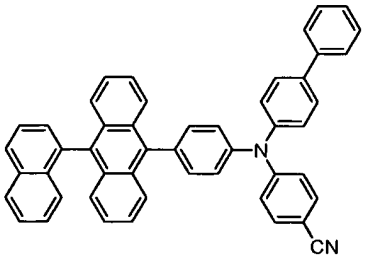

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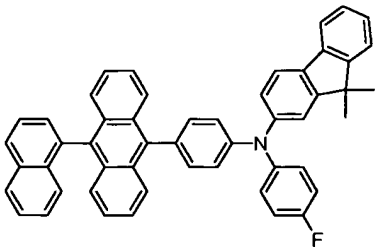

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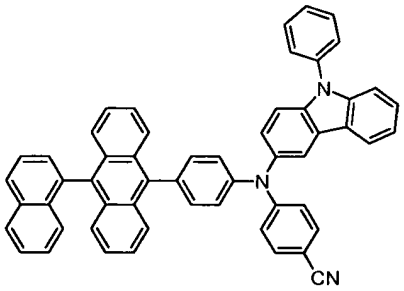

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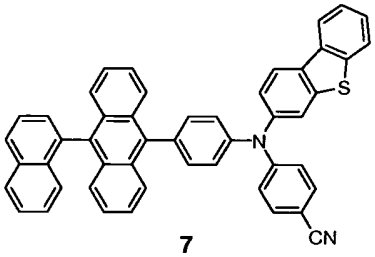

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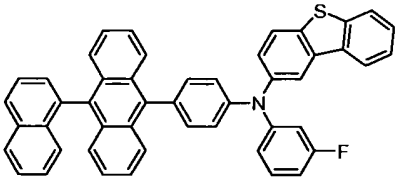

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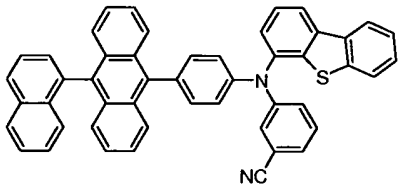

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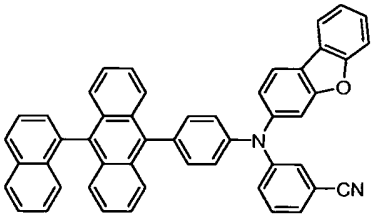

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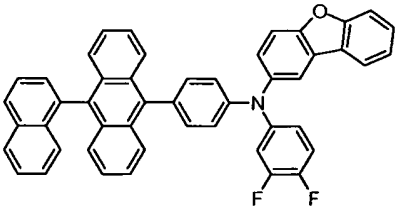

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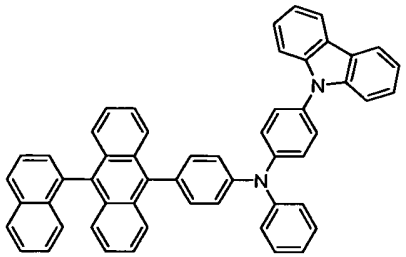

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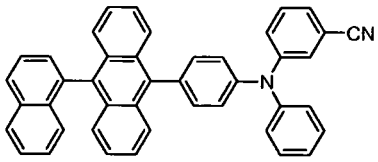

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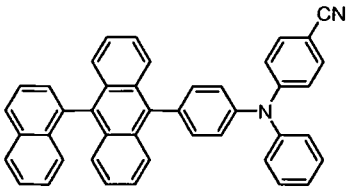

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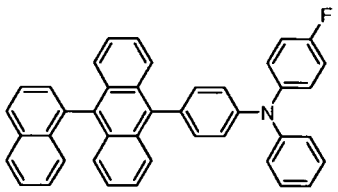

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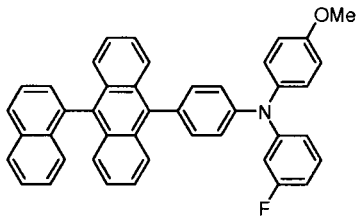

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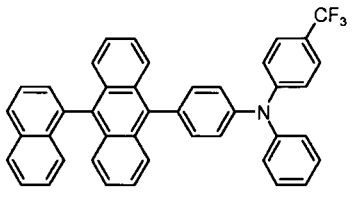

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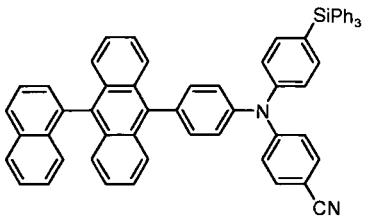

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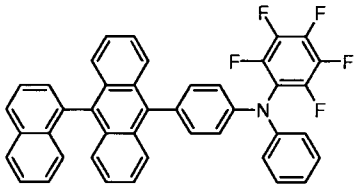

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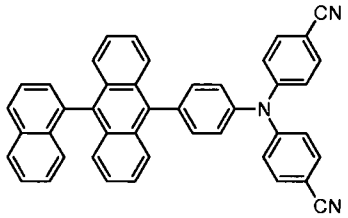

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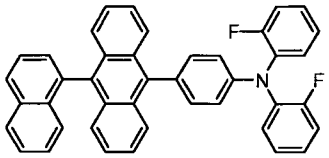

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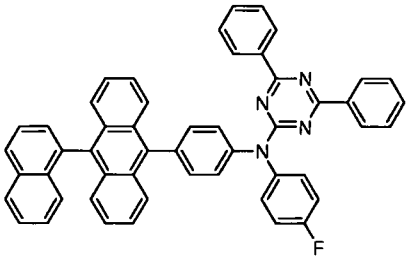

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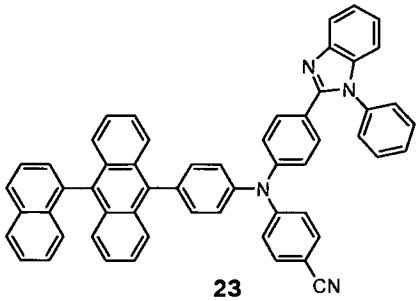

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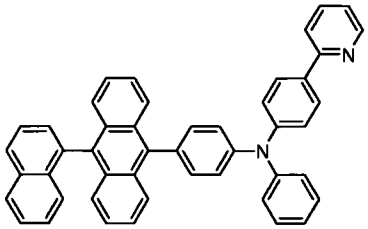

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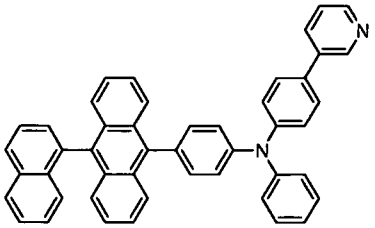

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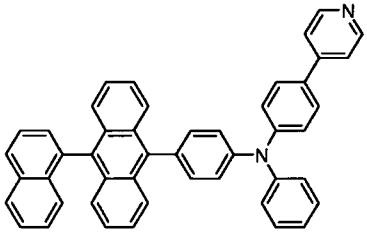

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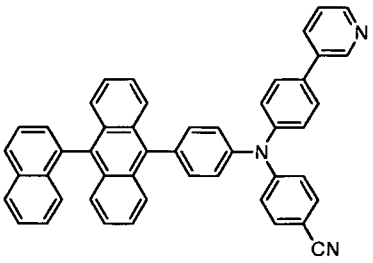

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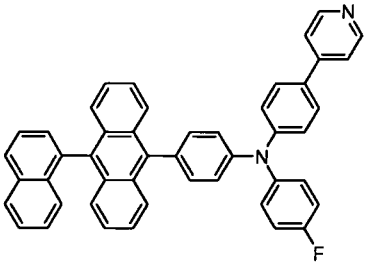

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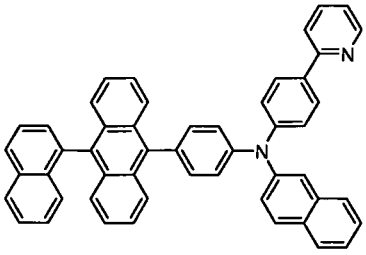

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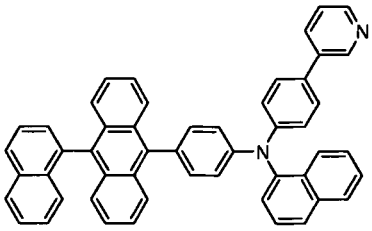

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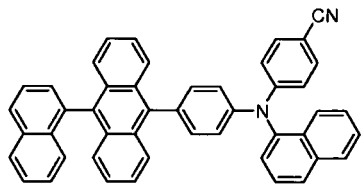

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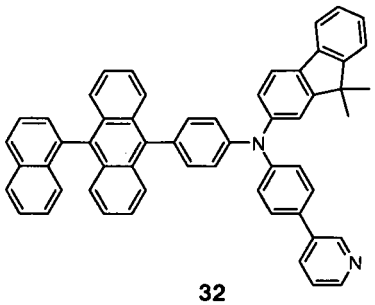

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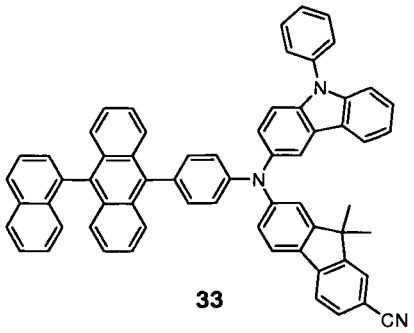

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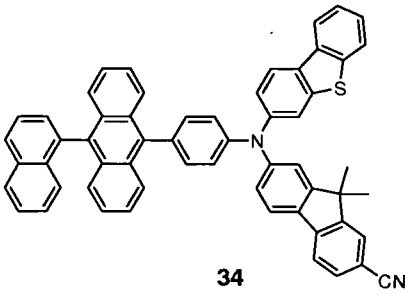

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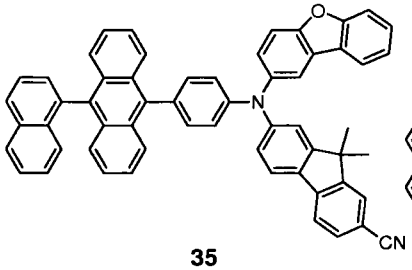

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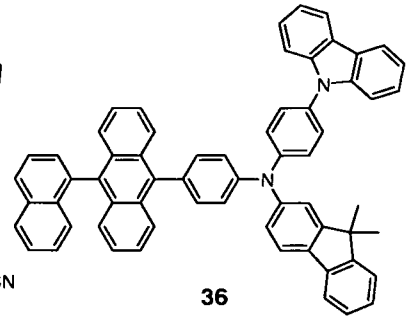

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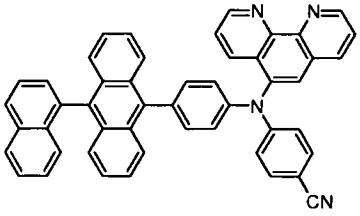

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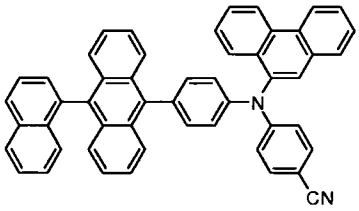

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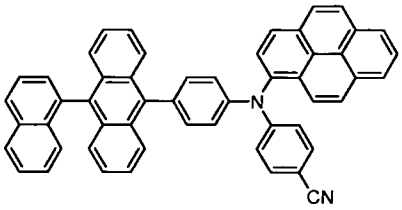

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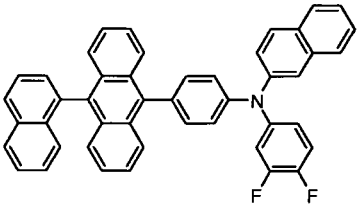

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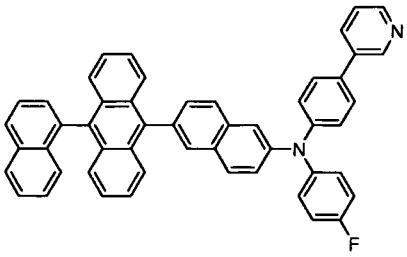

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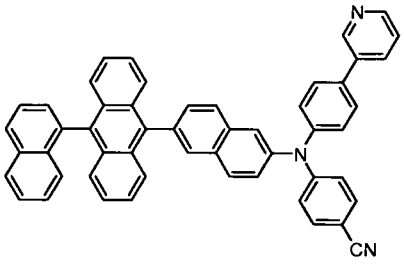

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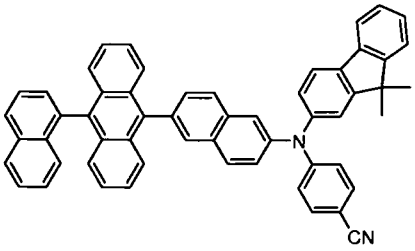

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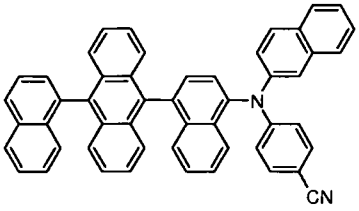

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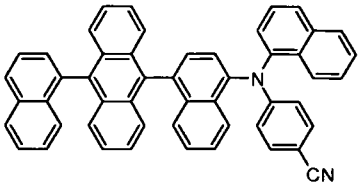

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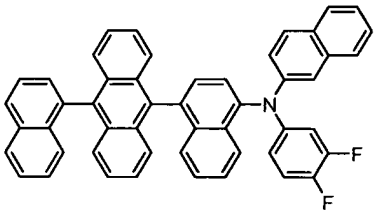

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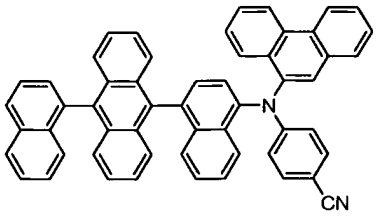

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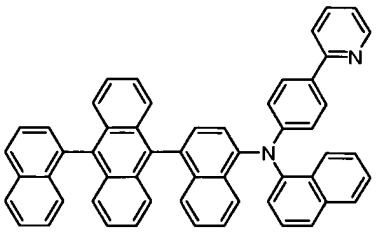

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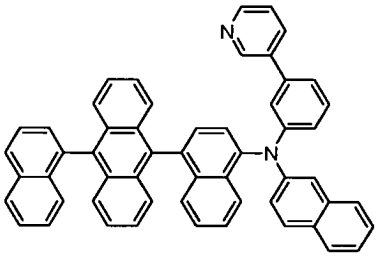

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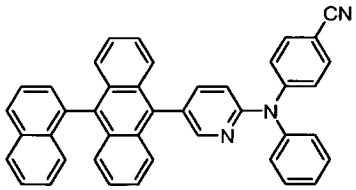

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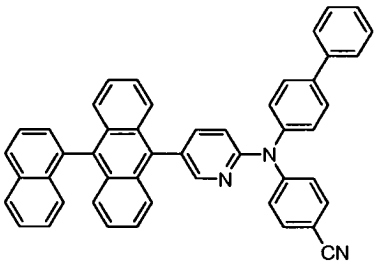

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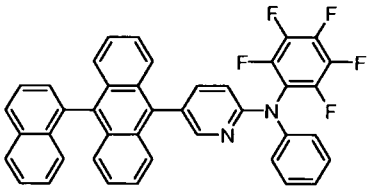

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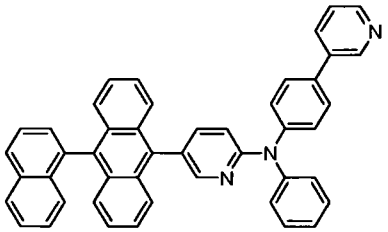

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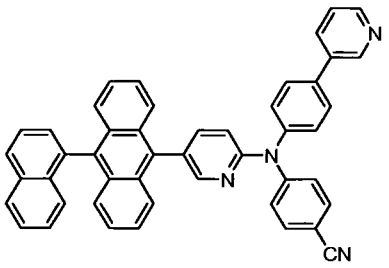

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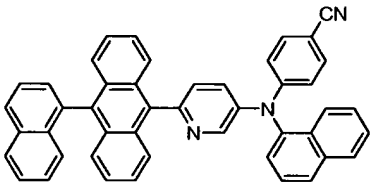

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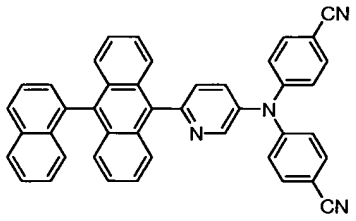

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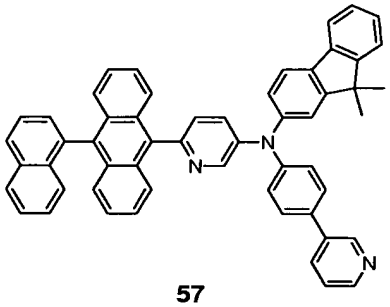

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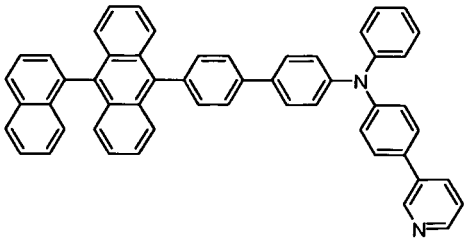

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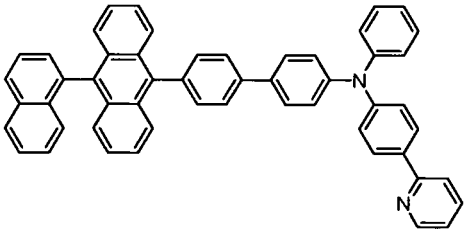

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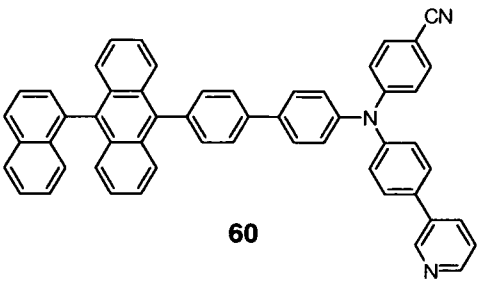

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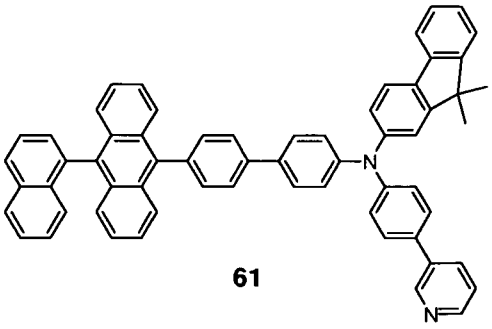

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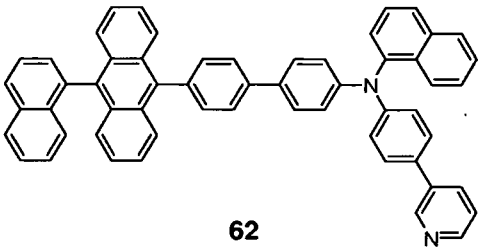

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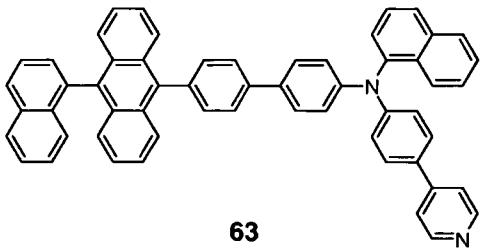

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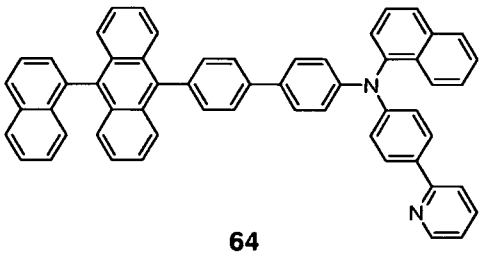

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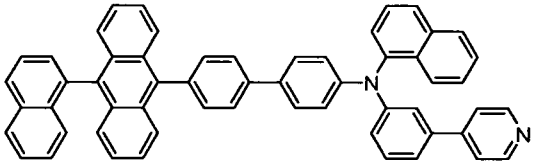

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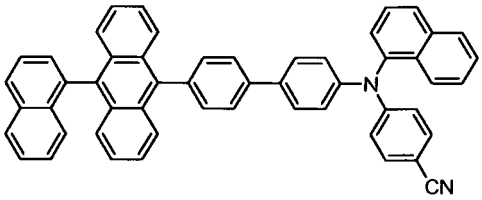

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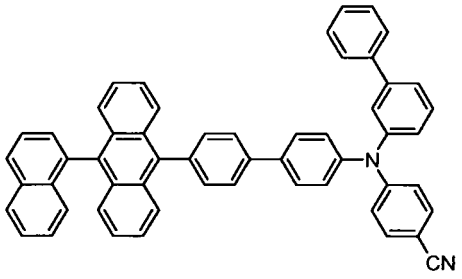

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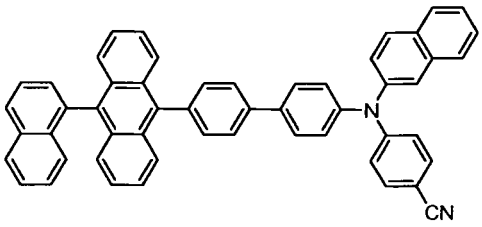

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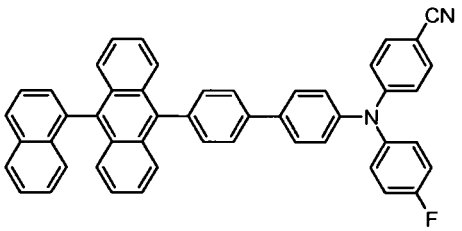

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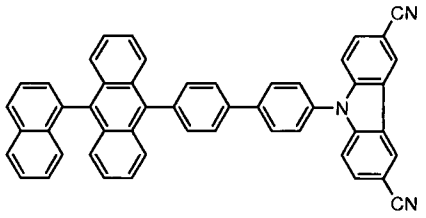

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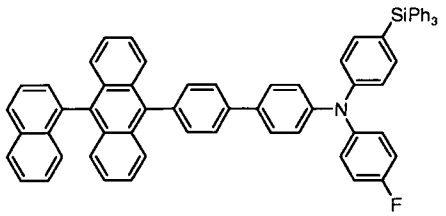

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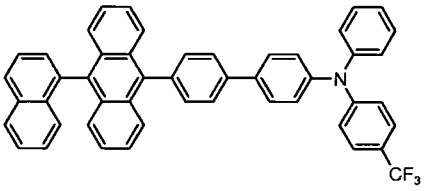

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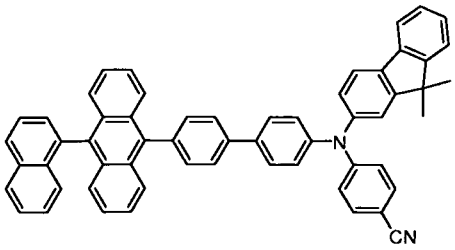

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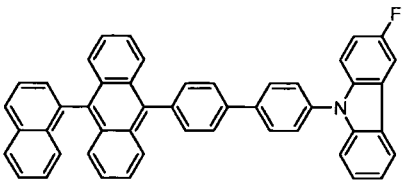

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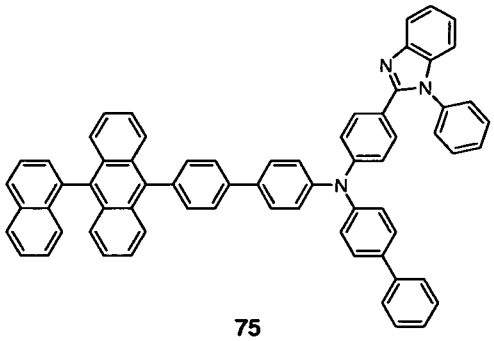

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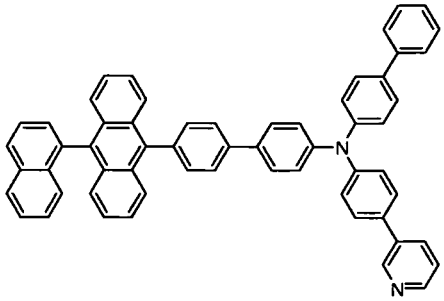

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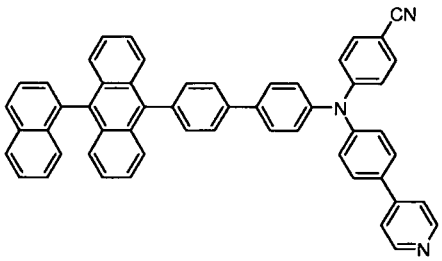

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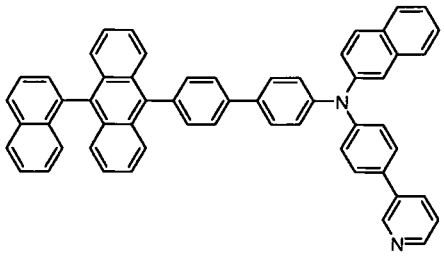

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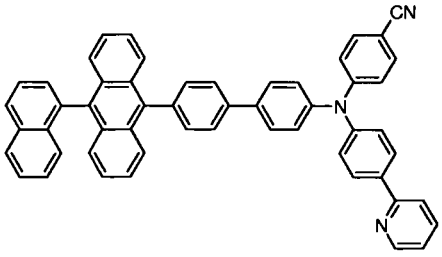

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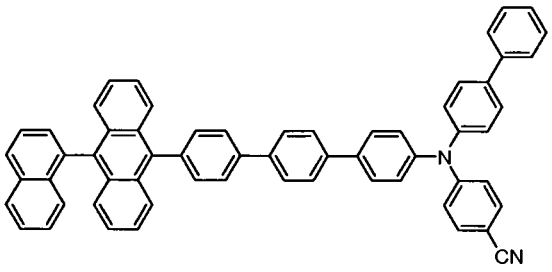

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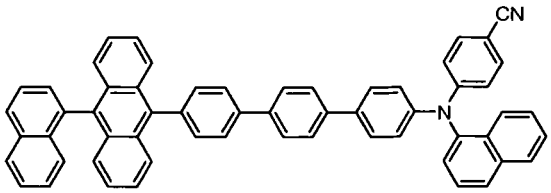

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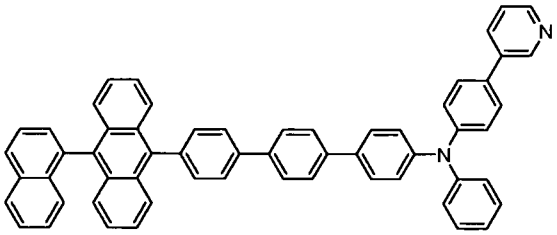

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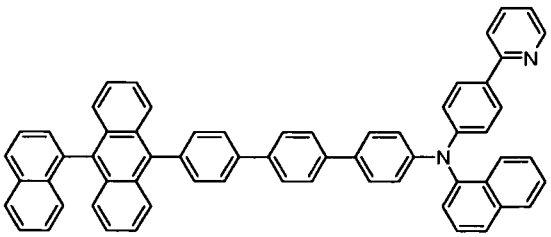

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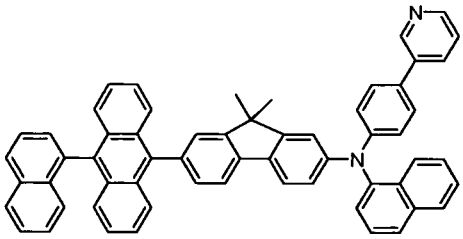

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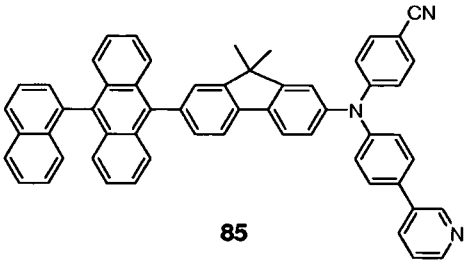

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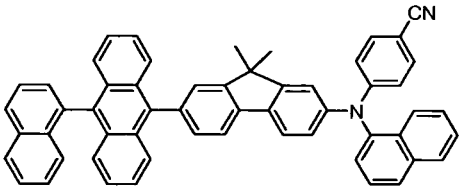

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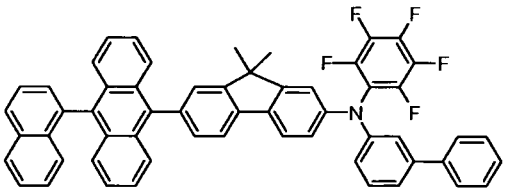

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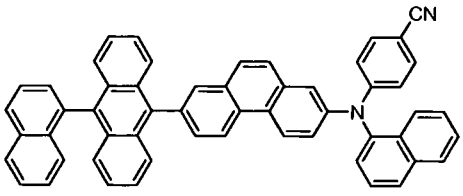

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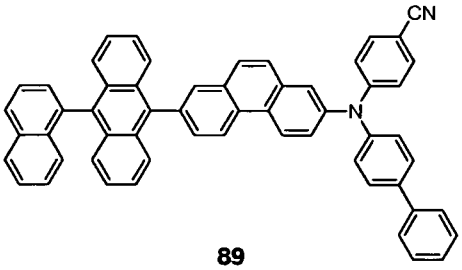

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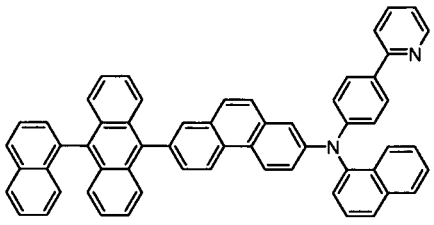

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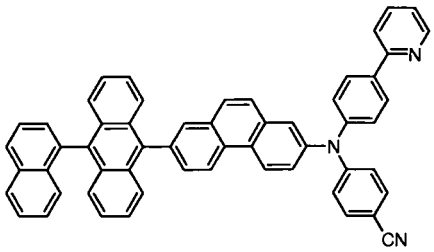

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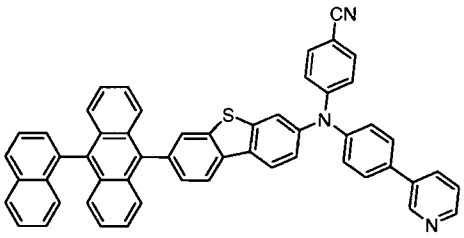

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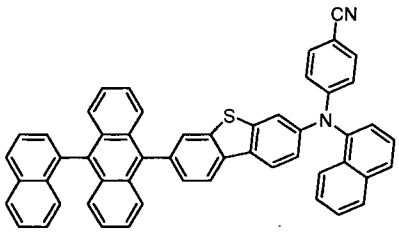

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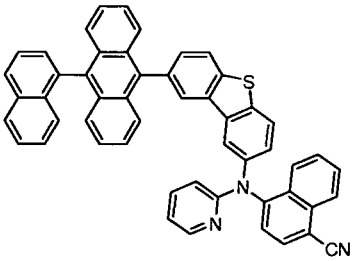

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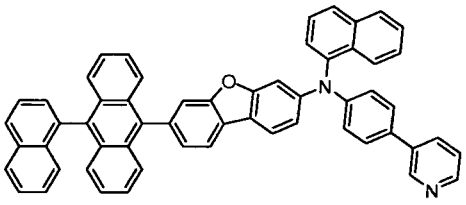

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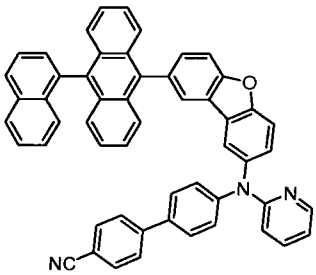

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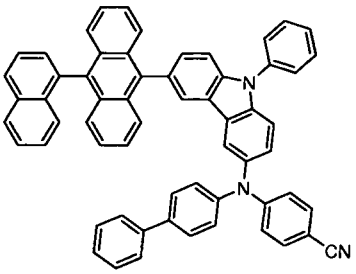

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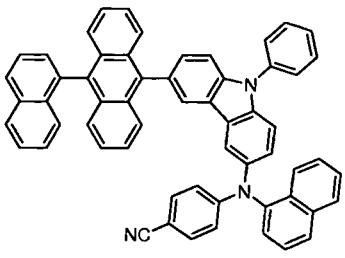

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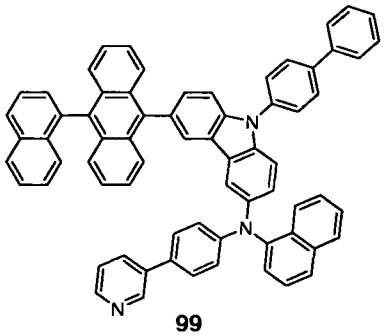

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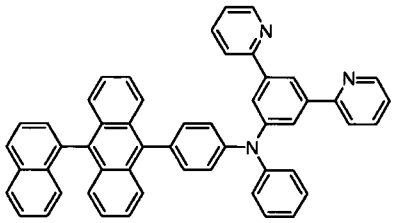

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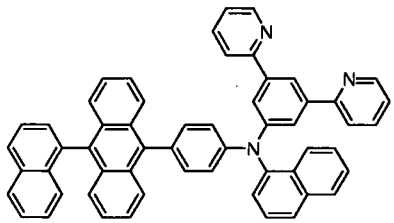

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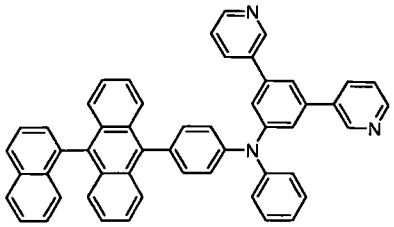

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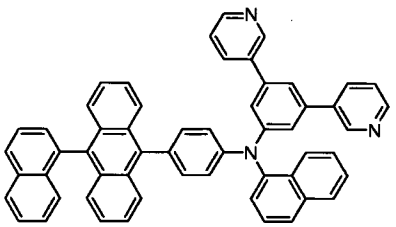

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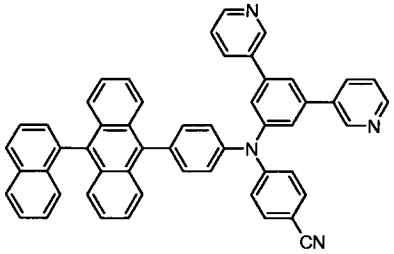

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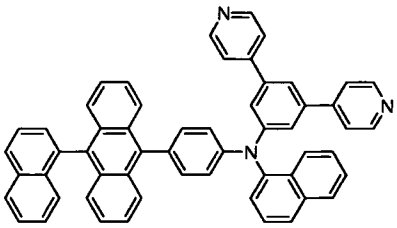

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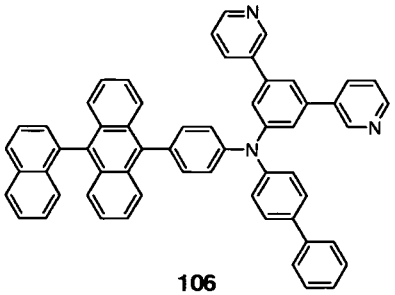

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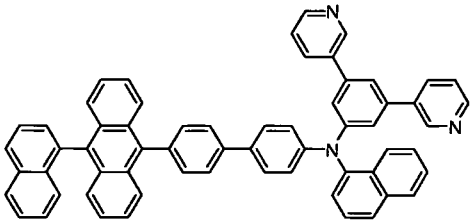

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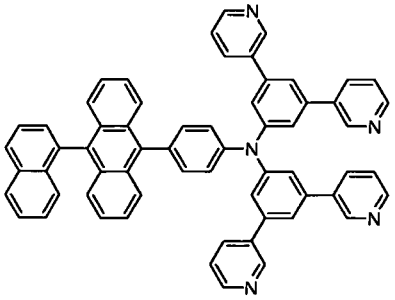

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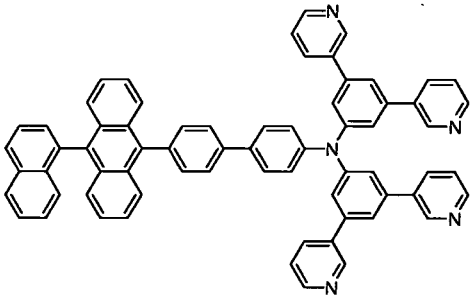
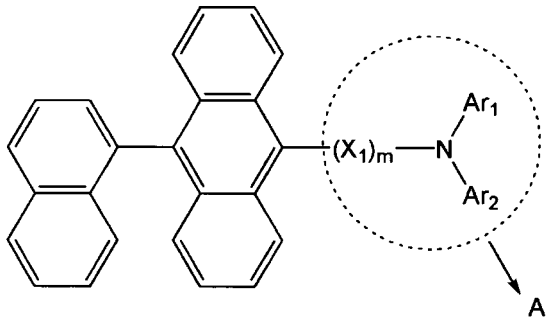


109

[0055] Since at least one of Ar_1 and Ar_2 in the amine-based compound of Formula 1 above is a C_6-C_{60} aryl group substituted with at least one electron withdrawing group selected from the group consisting of -F; -CN; -NO_2; a C_1-C_{60} alkyl group substituted with at least one -F; a C_2-C_{60} heteroaryl group; and a C_2-C_{60} heteroaryl group substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, -NO_2, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group, and a C_2-C_{60} heteroaryl group, a moiety represented by "A" in Formula 1' below may be able to withdraw electrons.

Formula 1'



[0056] The amine-based compound of Formula 1 has a naphthyl-anthracene core abundant in electrons and a moiety A with electron withdrawing ability, and thus the amine-based compound may have good electron transport characteristics. When the electron withdrawing group is a substituted or unsubstituted C_2-C_{60} heteroaryl group, the substituted or unsubstituted C_2-C_{60} heteroaryl group is linked to N of Formula 1, not directly, but via a C_6-C_{60} aryl group. Therefore, an organic light-emitting diode including the amine-based compound of Formula 1 may have improved efficiency characteristics. When the electron withdrawing group is -CN, an organic light-emitting diode including the amine-based compound of Formula 1 may have improved lifetime characteristics.

[0057] Not wishing to be bound by a particular theory, in either of i) an amine-based compound including a naphthyl-anthracene core that lacks the above-described electron withdrawing group or ii) an amine-based compound including a naphthyl-anthracene core with pyridine directly linked to N, the highest occupied molecular orbital (HOMO) electron density may be focused near the anthracene moiety. However, in the amine-based compound of Formula 1 above, the HOMO electron density may be dispersed near the amine moiety, so that the lowest unoccupied molecular orbital (LUMO) electron density may be relatively fixed near anthracene. This may result in the amine-based compound of Formula 1 having improved dipole characteristics. Thus, the electron transport characteristics of the amine-based compound of Formula 1 may be improved.

[0058] Therefore, an organic light-emitting diode including any of the amine-based compounds represented by Formula 1 above may have a low driving voltage, a high luminance, a high efficiency, and a long lifetime.

[0059] The amine-based compound of Formula 1 may be synthesized by any known organic synthesis method. A synthesis method of the amine-based compound of Formula 1 may be understood by those of ordinary skill in the art from the examples that will be described below.

[0060] At least one of the amine-based compounds according to the invention in its first aspect may be used between a pair of electrodes of an organic light-emitting diode. For example, at least one of the amine-based compounds may be in an EML and/or between a cathode and the EML (for example, in the ETL, the EIL, or a functional layer having both electron transport and electron injection capabilities).

[0061] According to another aspect of the present invention, an organic light-emitting diode includes a first electrode, a second electrode disposed opposite to the first electrode, and an organic layer disposed between the first electrode and the second electrode, wherein the organic layer includes at least one of the amine-based compounds of Formula 1 according to the invention in its first aspect.

[0062] As used herein, "(for example, the organic layer) including at least one amine-based compound" means (the organic layer) including one of the amine-based compounds of Formula 1 according to the invention in its first aspect, or at least two different amine-based compounds of Formula 1 according to the invention in its first aspect.

[0063] The organic layer may include only Compound 1 as the amine-based compound. Compound 1 may be in the EML or ETL of the organic light-emitting diode. The organic layer may include Compounds 1 and 3 as the amine-based compound. Compounds 1 and 3 may be in the same layer (for example, in the ETL) or may be in different layers (for example, in the EML and ETL, respectively).

[0064] The organic layer may include at least one layer selected from among a hole injection layer (HIL), a hole transport layer (HTL), a functional layer having both hole injection and hole transport capabilities (hereinafter, "H-functional layer"), a buffer layer, an electron blocking layer (EBL), an emission layer (EML), a hole blocking layer (HBL), an electron transport layer (ETL), an electron injection layer (EIL), and a functional layer having both electron injection and electron transport capabilities (hereinafter, "E-functional layer").

[0065] The term "organic layer" as used herein refers to a single layer and/or a plurality of layers disposed between the first and second electrodes of the organic light-emitting diode.

[0066] The organic layer may include an EML, wherein at least one of the amine-based compounds may be included in the EML.

[0067] The amine-based compound in the EML may serve as a host. When the amine-based compound in the EML

serves as a host, the EML may further include a fluorescent dopant. The fluorescent dopant may be a blue fluorescent dopant. The amine-based compound in the EML may serve as a dopant. When the amine-based compound in the EML serves as a dopant, the amine-based compound may be a blue fluorescent dopant.

[0068] The organic layer may include an ETL, wherein at least one of the amine-based compounds may be included in the ETL.

[0069] Figure 1 is a schematic sectional view of an organic light-emitting diode 10 according to an embodiment of the invention. Hereinafter, a structure of an organic light-emitting diode according to an embodiment of the invention and a method of manufacturing the same will now be described with reference to Figure 1.

[0070] The substrate 11 may be any substrate that is used in existing organic light-emitting diodes. The substrate 11 may be a glass substrate or a transparent plastic substrate with strong mechanical strength, thermal stability, transparency, surface smoothness, ease of handling, and water resistance.

[0071] The first electrode 13 may be formed by depositing or sputtering a first electrode-forming material on the substrate 11. When the first electrode 13 constitutes an anode, a material having a high work function may be used as the first electrode-forming material to facilitate hole injection. The first electrode 13 may be a reflective electrode or a transmission electrode. Suitable first electrode-forming materials include transparent and conductive materials such as ITO, IZO, SnO_2, and ZnO. The first electrode 13 may be formed as a reflective electrode using magnesium (Mg), aluminium (Al), aluminium-lithium (Al-Li), calcium (Ca), magnesium-indium (Mg-In), magnesium-silver (Mg-Ag), or the like.

[0072] The first electrode 13 may have a single-layer structure or a multi-layer structure including at least two layers. For example, the first electrode 13 may have a three-layered structure of ITO/Ag/ITO, but is not limited thereto.

[0073] The organic layer 15 may be disposed on the first electrode 13.

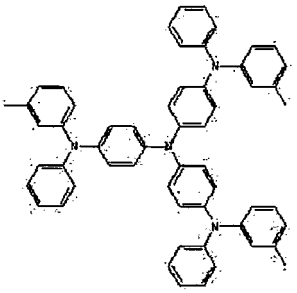
[0074] The organic layer 15 may include a HIL, a HTL, a buffer layer, an EML, an ETL, and an EIL.

[0075] The HIL may be formed on the first electrode 13 by vacuum deposition, spin coating, casting, Langmuir-Blodgett (LB) deposition, or the like.

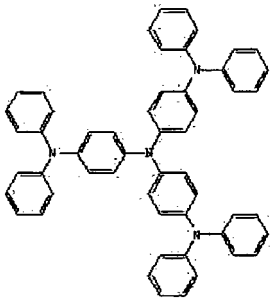
[0076] When the HIL is formed using vacuum deposition, vacuum deposition conditions may vary according to the compound that is used to form the HIL, and the desired structure and thermal properties of the HIL to be formed. For example, vacuum deposition may be performed at a temperature of about 100°C to about 500°C, a pressure of about 0.133 x 10^{-6} Pa (10^{-8} torr) to about 0.133 Pa (10^{-3} torr), and a deposition rate of about 0.001 to about 10 nm/sec (about 0.01 to about 100 Å/sec). However, the deposition conditions are not limited thereto.

[0077] When the HIL is formed using spin coating, the coating conditions may vary according to the compound that is used to form the HIL, and the desired structure and thermal properties of the HIL to be formed. For example, the coating rate may be in the range of about 2000 rpm to about 5000 rpm, and a temperature at which heat treatment is performed to remove a solvent after coating may be in the range of about 80°C to about 200°C. However, the coating conditions are not limited thereto.

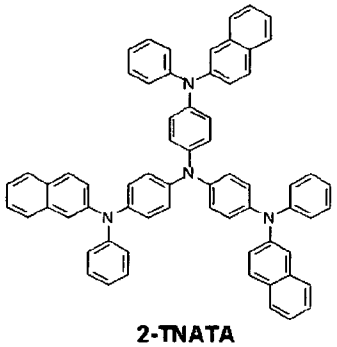
[0078] The HIL 130 may comprise any material that is commonly used to form a HIL. Examples of the material that can be used to form the HIL are, but are not limited to, N,N'-diphenyl-N,N'-bis-[4-(phenyl-m-tolyl-amino)-phenyl]-biphenyl-4,4'-diamine, (DNTPD), a phthalocyanine compound such as copper phthalocyanine, 4,4',4"-tris (3-methylphenylphenylamino)triphenylamine (m-MTDATA), N,N'-di(1-naphthyl)-N,N'-diphenylbenzidine (NPB), 4,4'4"-Tris(N,N-diphenylamino)triphenylamine (TDATA), 4,4'4"-Tris-(N-(naphthylen-2-yl)-N-phenylamine)triphenylamine (2-TNATA), polyaniline/dodecylbenzenesulfonic acid (Pani/DBSA), poly(3,4-ethylenedioxythiophene)/poly(4-styrenesulfonate) (PEDOT/PSS), polyaniline/camphor sulfonic acid (Pani/CSA), and polyaniline/poly(4-styrenesulfonate) (PANI/PSS).



m-MTDATA



TDATA

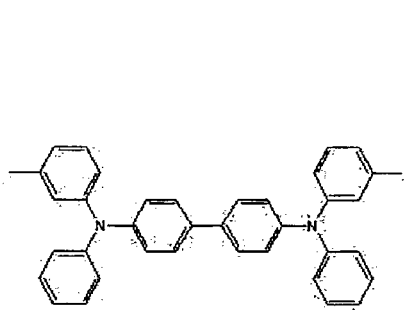


2-TNATA

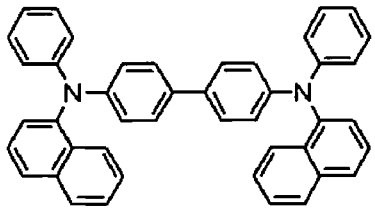
[0079] The thickness of the HIL may be about 10 nm (100 Å) to about 1000 nm (10000 A), and preferably may be about 10 nm (100 Å) to about 100 nm (1000 Å). When the thickness of the HIL is within these ranges, the HIL may have good hole injecting ability without a substantial increase in driving voltage.

[0080] Then, a HTL may be formed on the HIL by using vacuum deposition, spin coating, casting, Langmuir-Blodgett (LB) deposition, or the like. When the HTL is formed using vacuum deposition or spin coating, the conditions for deposition and coating may be similar to those for the formation of the HIL, though the conditions for the deposition and coating may vary according to the material that is used to form the HTL.

[0081] Non-limiting examples of suitable known hole transport materials are carbazole derivatives, such as N-phenylcarbazole or polyvinylcarbazole, N,N'-bis(3-methylphenyl)-N,N'-diphenyl-[1,1-biphenyl]-4,4'-diamine (TPD), 4,4',4"-tris(N-carbazolyl)triphenylamine (TCTA), and N,N'-di(1-naphthyl)-N,N'-diphenylbenzidine) (NPB).



TPD



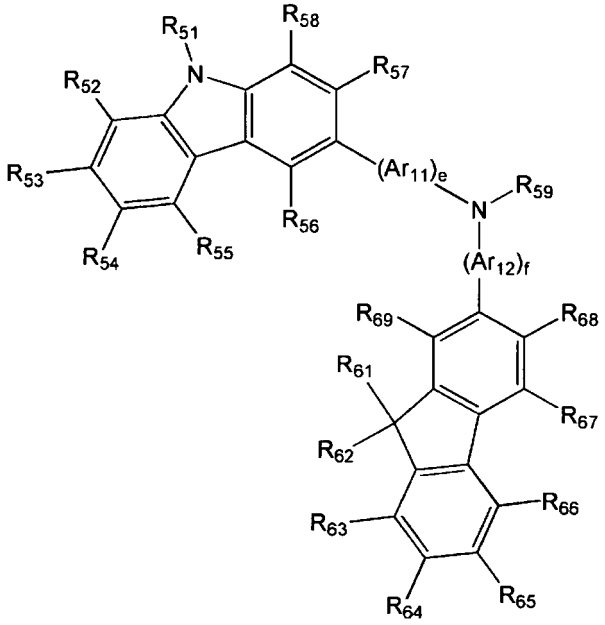
NPB

[0082] The thickness of the HTL may be about 5 nm (50 Å) to about 200 nm (2000 Å), and preferably may be about 10 nm (100 Å) to about 150 nm (1500 Å). When the thickness of the HTL is within these ranges, the HTL may have good hole transporting ability without a substantial increase in driving voltage.

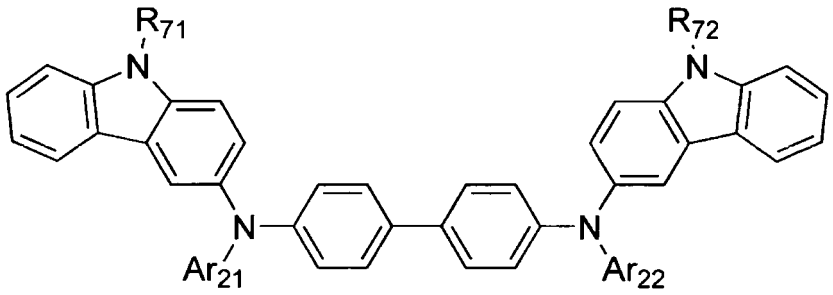
[0083] The H-functional layer (having both hole injection and hole transport capabilities) may contain at least one material from each group of the hole injection layer materials and hole transport layer materials. The thickness of the H-functional layer may be from about 10 nm (100Å) to about 1000 nm (10,000Å), and preferably may be from about 10 nm (100Å) to about 100 nm (1,000Å). When the thickness of the H-functional layer is within these ranges, the H-functional layer may have good hole injection and transport capabilities without a substantial increase in driving voltage.

[0084] At least one of the HIL, HTL, and H-functional layer may include at least one of a compound of Formula 300 below and a compound of Formula 350 below:

Formula 300



Formula 350



[0085] In each of Formulae 300 and 350, Ar_{11} and Ar_{12} may each independently be a substituted or unsubstituted C_6-C_{60} arylene group and Ar_{21} and Ar_{22} may each independently be a substituted or unsubstituted C_6-C_{60} aryl group. Ar_{11} and Ar_{12} may each independently be one of a phenylene group, a naphthylene group, a phenanthrenylene group, and a pyrenylene group; and a phenylene group, a naphthylene group, a phenanthrenylene group, a fluorenylene group, and a pyrenylene group that are substituted with at least one of a phenyl group, a naphthyl group, and an anthryl group. Ar_{21} and Ar_{22} may each independently be one of a phenyl group, a naphthyl group, a phenanthrenyl group, and a pyrenyl group; and a phenyl group, a naphthyl group, a phenanthrenyl group, a fluorenyl group, and a pyrenyl group that are substituted with at least one of a phenyl group, a naphthyl group, and an anthryl group.

[0086] In Formula 300, e and f may each independently be an integer from 0 to 5, for example, may be 0, 1, or 2. In a non-limiting embodiment, e may be 1, and f may be o.

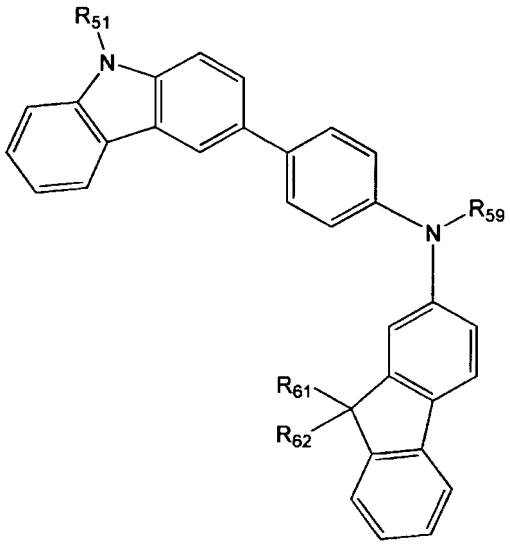
[0087] In each of Formulae 300 and 350, R_{51} to R_{58}, R_{61} to R_{69}, R_{71}, and R_{72} may each independently be a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a substituted or unsubstituted C_1-C_{60} alkyl group, a substituted or unsubstituted C_2-C_{60} alkenyl group, a substituted or unsubstituted C_2-C_{60} alkynyl group, a substituted or unsubstituted C_1-C_{60} alkoxy group, a substituted or unsubstituted C_3-C_{60} cycloalkyl group, a substituted or unsubstituted C_6-C_{60a} aryl group, a substituted or unsubstituted C_6-C_{60} aryloxy group, or a substituted or unsubstituted C_6-C_{60} arylthiol group. In some non-limiting embodiments, R_5, to R_{58}, R_{61} to R_{69}, R_{71}, and R_{72} may each independently be one of a hydrogen atom; a deuterium atom; a halogen atom; a hydroxyl group; a cyano group; $-NO_2$; an amino group; an amidino group; hydrazine; hydrazone; a carboxyl group or a salt thereof; a sulfonic acid group or a salt thereof; a phosphoric acid or a salt thereof; a C_1-C_{10} alkyl group (for example, a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, a hexyl group, or the like); a C_1-C_{10} alkoxy group (for example, a methoxy group, an ethoxy group, a propoxy group, a butoxy group, a pentoxy group, or the like); a C_1-C_{10} alkyl group and a C_1-C_{10} alkoxy group that are substituted with at least one of a deuterium

atom, a halogen atom, a hydroxyl group, a cyano group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, and a phosphoric acid or a salt thereof; a phenyl group; a naphthyl group; an anthryl group; a fluorenyl group; a pyrenyl group; and a phenyl group, a naphthyl group, an anthryl group, a fluorenyl group, and a pyrenyl group that are substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{10} alkyl group, and a C_1-C_{10} alkoxy group.

[0088] In Formula 300, R_{59} may be one of a phenyl group; a naphthyl group; an anthryl group; a biphenyl group; a pyridyl group; and a phenyl group, a naphthyl group, an anthryl group, a biphenyl group, and a pyridyl group that are substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a substituted or unsubstituted C_1-C_{20} alkyl group, and a substituted or unsubstituted C_1-C_{20} alkoxy group.

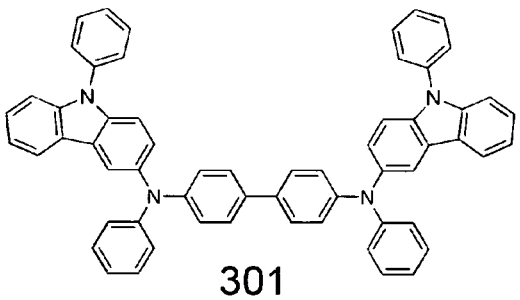
[0089] In an embodiment the compound of Formula 300 is a compound represented by Formula 300A below:

Formula 300A



[0090] R_{51}, R_{60}, R_{61} and R_{59} in Formula 300A are as defined above, and thus a detailed description thereof will not be provided herein.

[0091] At least one of the HIL, HTL, and H-functional layer may include at least one of compounds represented by Formulae 301 to 320 below:



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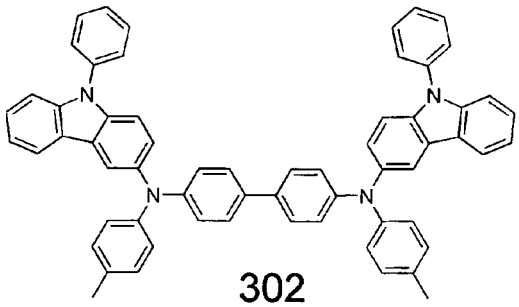

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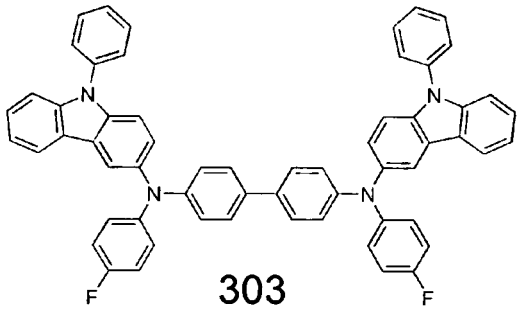
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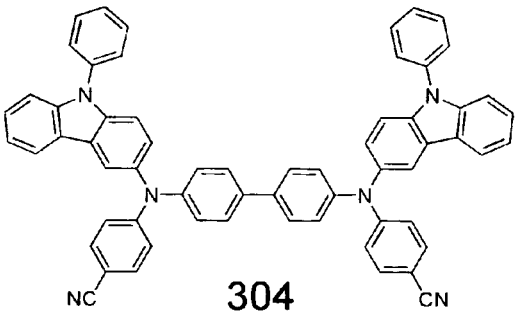
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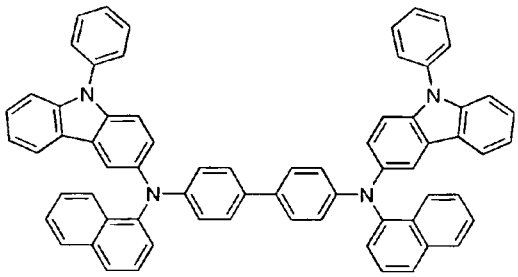
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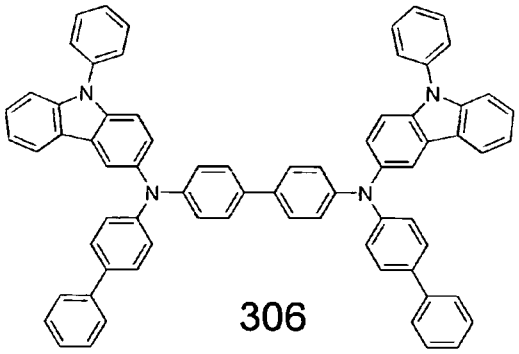
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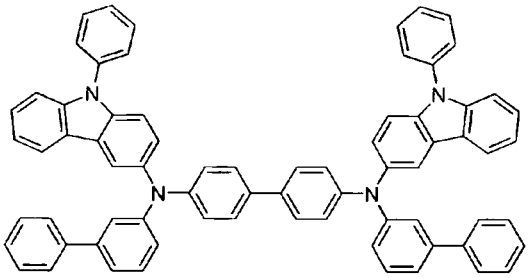
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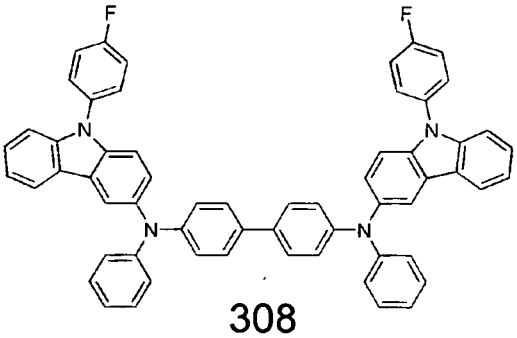
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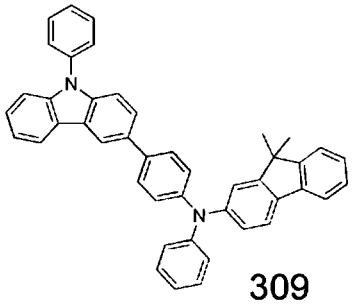
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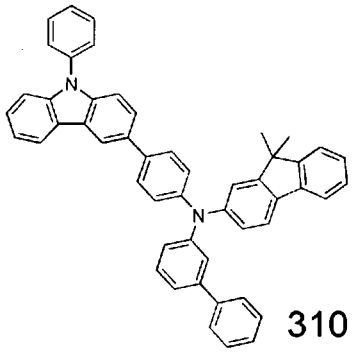
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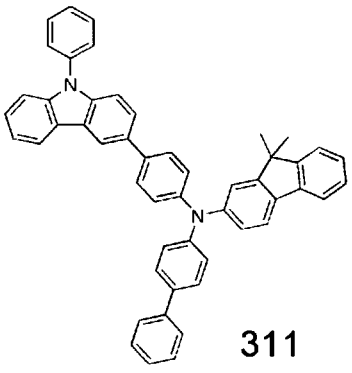
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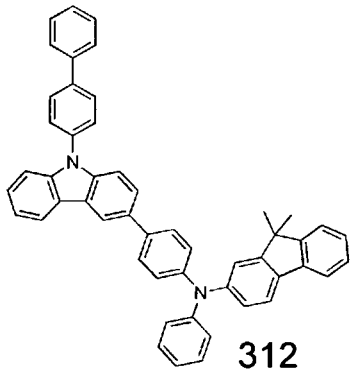
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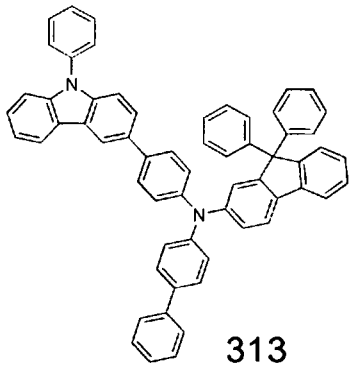
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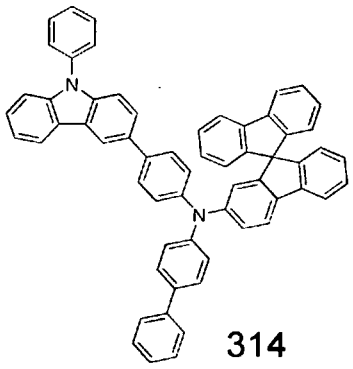
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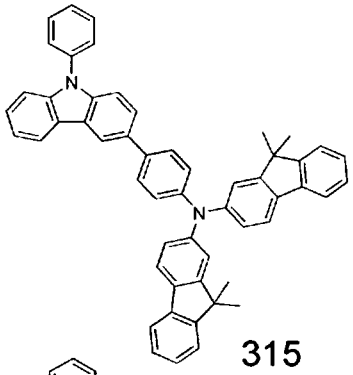
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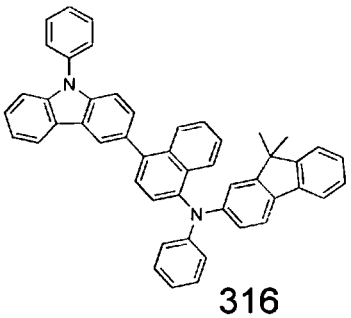
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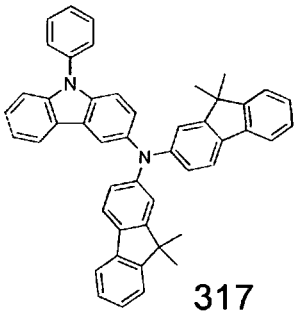
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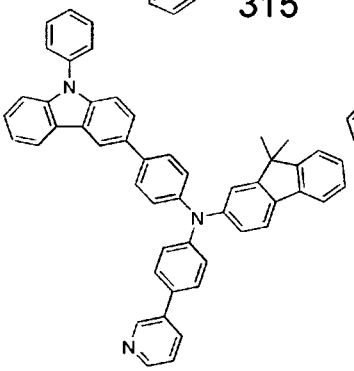
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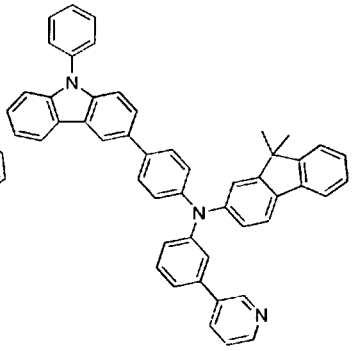
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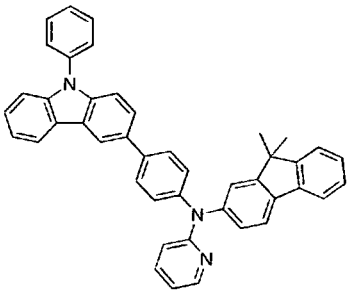
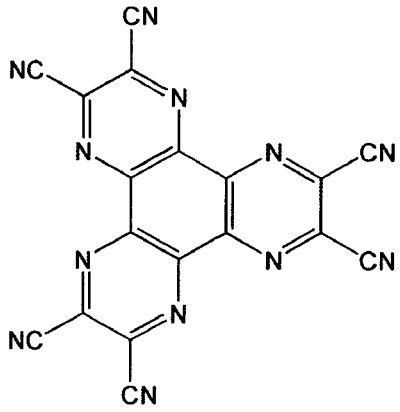

320

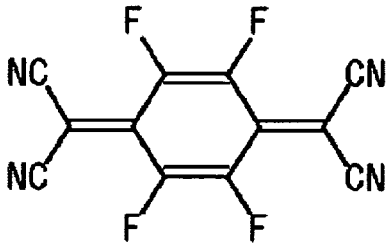
[0092] At least one of the HIL, HTL, and H-functional layer may further include a charge-generating material for improved layer conductivity, in addition to a known hole injecting material, hole transport material, and/or material having both hole injection and hole transport capabilities as described above.

[0093] The charge-generating material may be, for example, a p-dopant. The p-dopant may be one of quinine derivatives, metal oxides, and compounds with a cyano group, but are not limited thereto. Non-limiting examples of the p-dopant include quinone derivatives such as tetracyanoquinonedimethane (TCNQ), 2,3,5,6-tetrafluoro-tetracyano-1,4-benzoquinonedimethane (F_4-TCNQ), and the like; metal oxides such as tungsten oxide, molybdenum oxide, and the like; and cyano-containing compounds such as Compound 200 below.

Compound 200



F4-TCNQ



[0094] When the hole injection layer, hole transport layer, or H-functional layer further includes a charge-generating material, the charge-generating material may be homogeneously dispersed or inhomogeneously distributed in the layer.

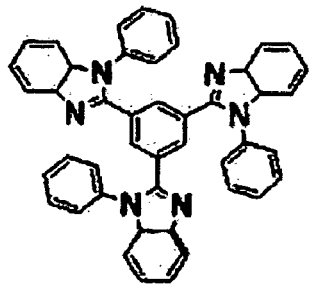
[0095] A buffer layer may be disposed between at least one of the HIL, HTL, and H-functional layer, and the EML. The buffer layer may compensate for an optical resonance distance of light according to a wavelength of the light emitted from the EML, and thus may increase efficiency. The buffer layer may include any hole injecting material or hole transporting material that is widely known. The buffer layer may include the same material as one of the materials included in the HIL, HTL, and H-functional layer that lies beneath the buffer layer.

[0096] Then, an EML may be formed on the HTL, H-functional layer, or buffer layer by vacuum deposition, spin coating, casting, Langmuir-Blodget (LB) deposition, or the like. When the EML is formed using vacuum deposition or spin coating, the deposition and coating conditions may be similar to those for the formation of the HIL, though the conditions for deposition and coating may vary according to the material that is used to form the EML.

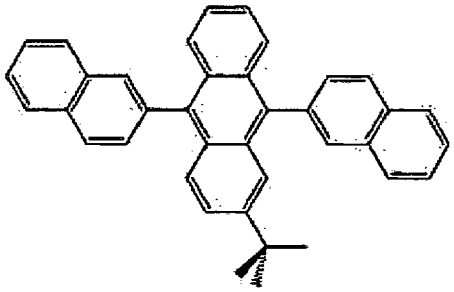
[0097] The EML may include at least one of the amine-based compounds of Formula 1.

[0098] The amine-based compound in the EML may serve as a dopant, for example, as a blue fluorescent dopant. The EML may further include a host, in addition to the amine-based compound.

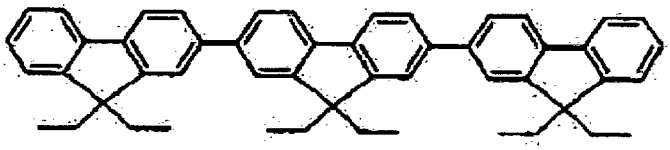
[0099] Example of the host are Alq_3, 4,4'-N,N'-dicarbazole-biphenyl (CBP), poly(n-vinylcarbazole) (PVK), 9,10-di(naphthalene-2-yl)anthracene (ADN), tris(4-carbazoyl-9-ylphenyl)amine (TCTA), 1,3,5-tris(N-phenylbenzimidazole-2-yl)benzene (TPBI), 3-tert-butyl-9,10-di-2-naphthylanthracene (TBADN), E3, distyrylarylene (DSA), dmCBP (see formula below), and Compounds 501 to 509 below, but are not limited thereto.



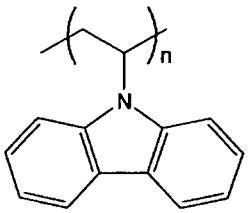
TPBI



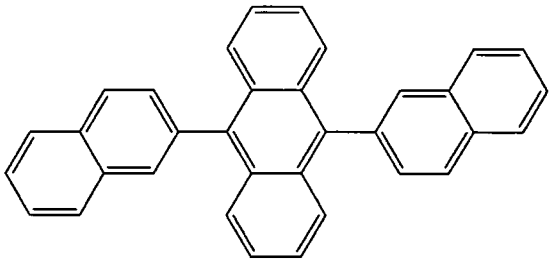
TBADN



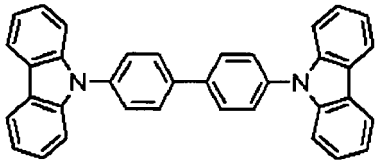
E3



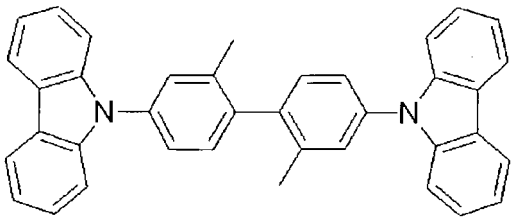
PVK



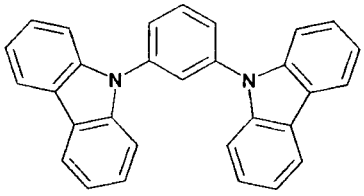
ADN



CBP



dmCBP



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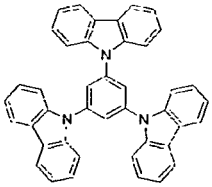

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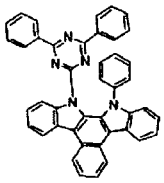

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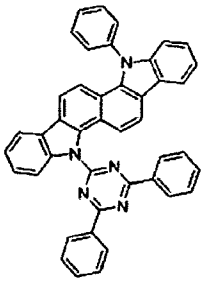

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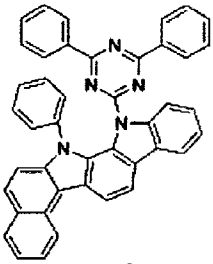

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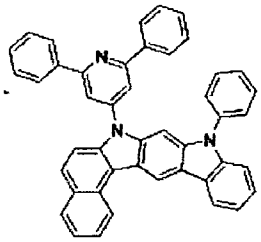

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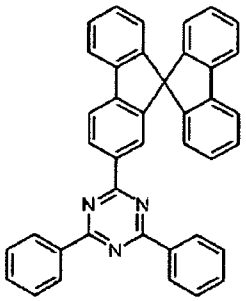

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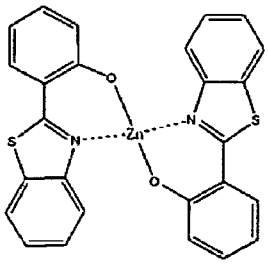

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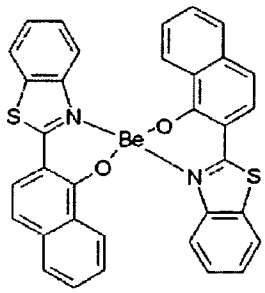
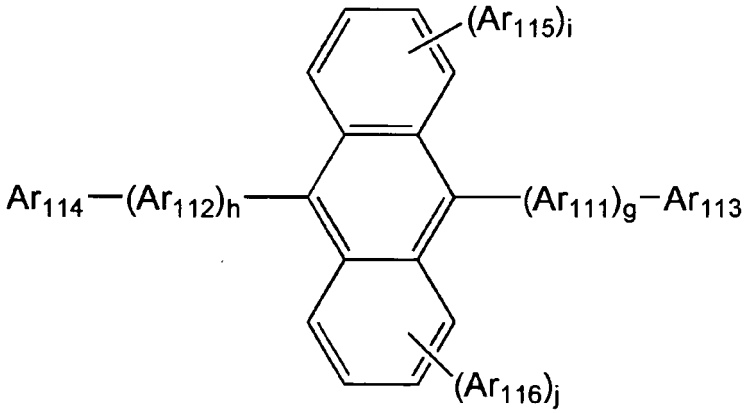


509

[0100] An anthracene-based compound represented by Formula 400 below may be used as the host.

Formula 400

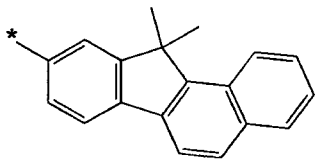


[0101] In Formula 400, Arm and Ar_{112} may each independently be a substituted or unsubstituted C_6-C_{60} arylene group; Ar_{113} to Ar_{116} may each independently be a substituted or unsubstituted C_1-C_{10} alkyl group or a substituted or unsubstituted C_6-C_{60} aryl group; and g, h, i, and j may each independently be an integer selected from 0, 1, 2, 3 and 4.

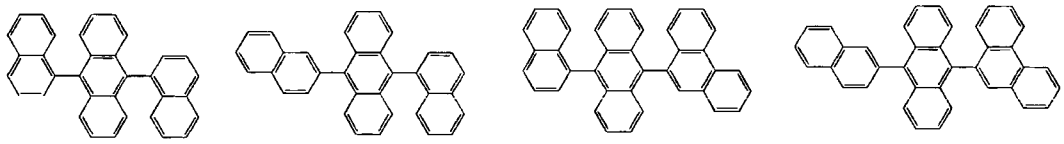
[0102] Arm and Ar_{112} in Formula 400 may each independently be a phenylene group, a naphthylene group, a phenanthrenylene group, or a pyrenylene group; or a phenylene group, a naphthylene group, a phenanthrenylene group, a fluorenyl group, or a pyrenylene group that are substituted with at least one of a phenyl group, a naphthyl group, and an anthryl group.

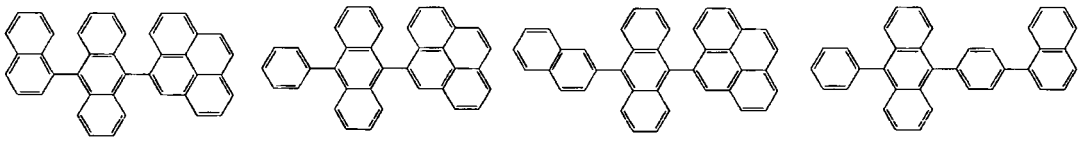
[0103] In Formula 400 above, g, h, i, and j may each independently be 0, 1, or 2.

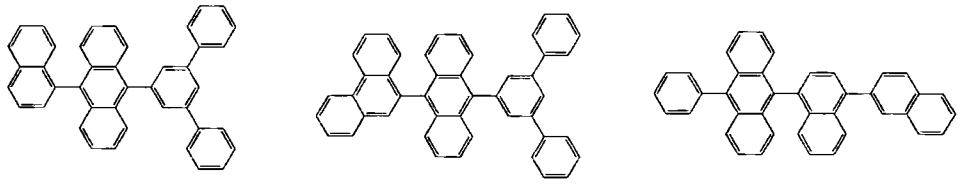
[0104] Ar_{113} to Ar_{116} in Formula 400 may each independently be one of a C_1-C_{10} alkyl group substituted with at least one of a phenyl group, a naphthyl group, and an anthryl group; a phenyl group; a naphthyl group; an anthryl group; a pyrenyl group; a phenanthrenyl group; a fluorenyl group; a phenyl group, a naphthyl group, an anthryl group, a pyrenyl group, a phenanthrenyl group, and a fluorenyl group that are substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_1-C_{60} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a pyrenyl group, a phenanthrenyl group, and a fluorenyl group; and

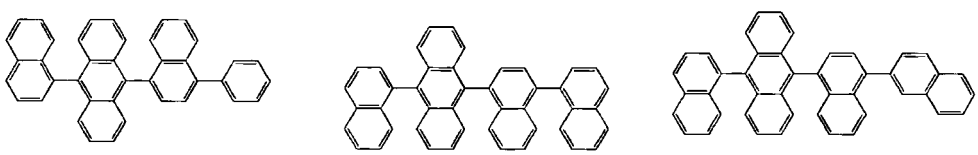


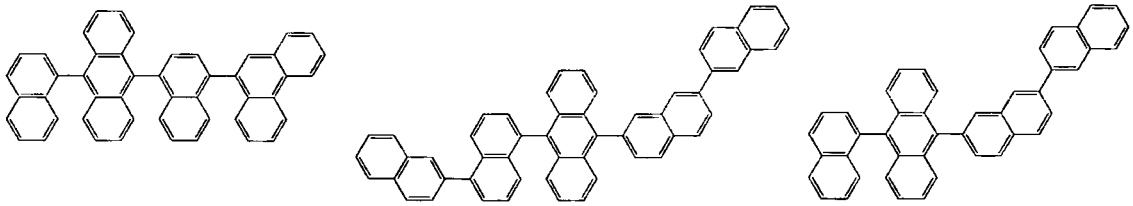
[0105] For example, the anthracene compound of Formula 400 above may be one of the compounds represented by the following formulae, but is not limited thereto:

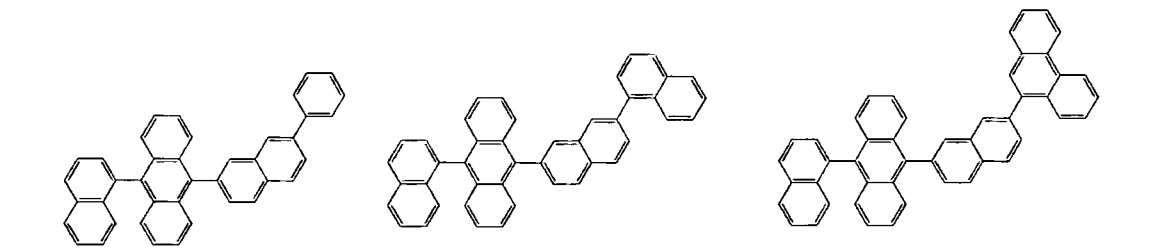


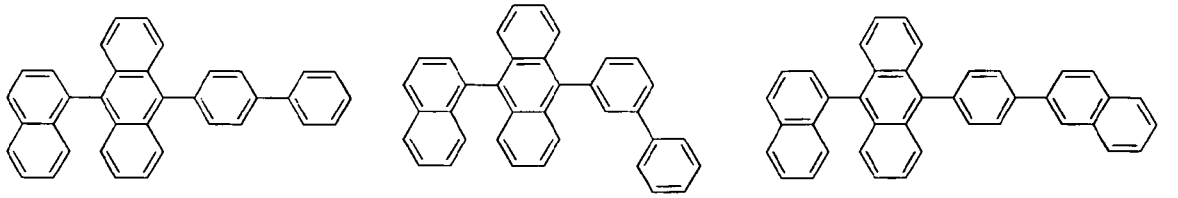


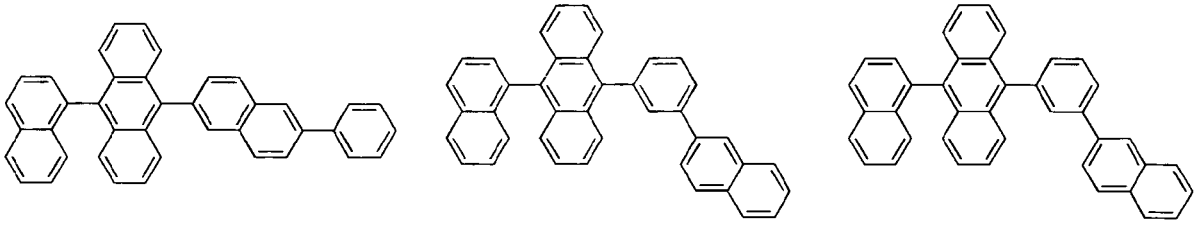


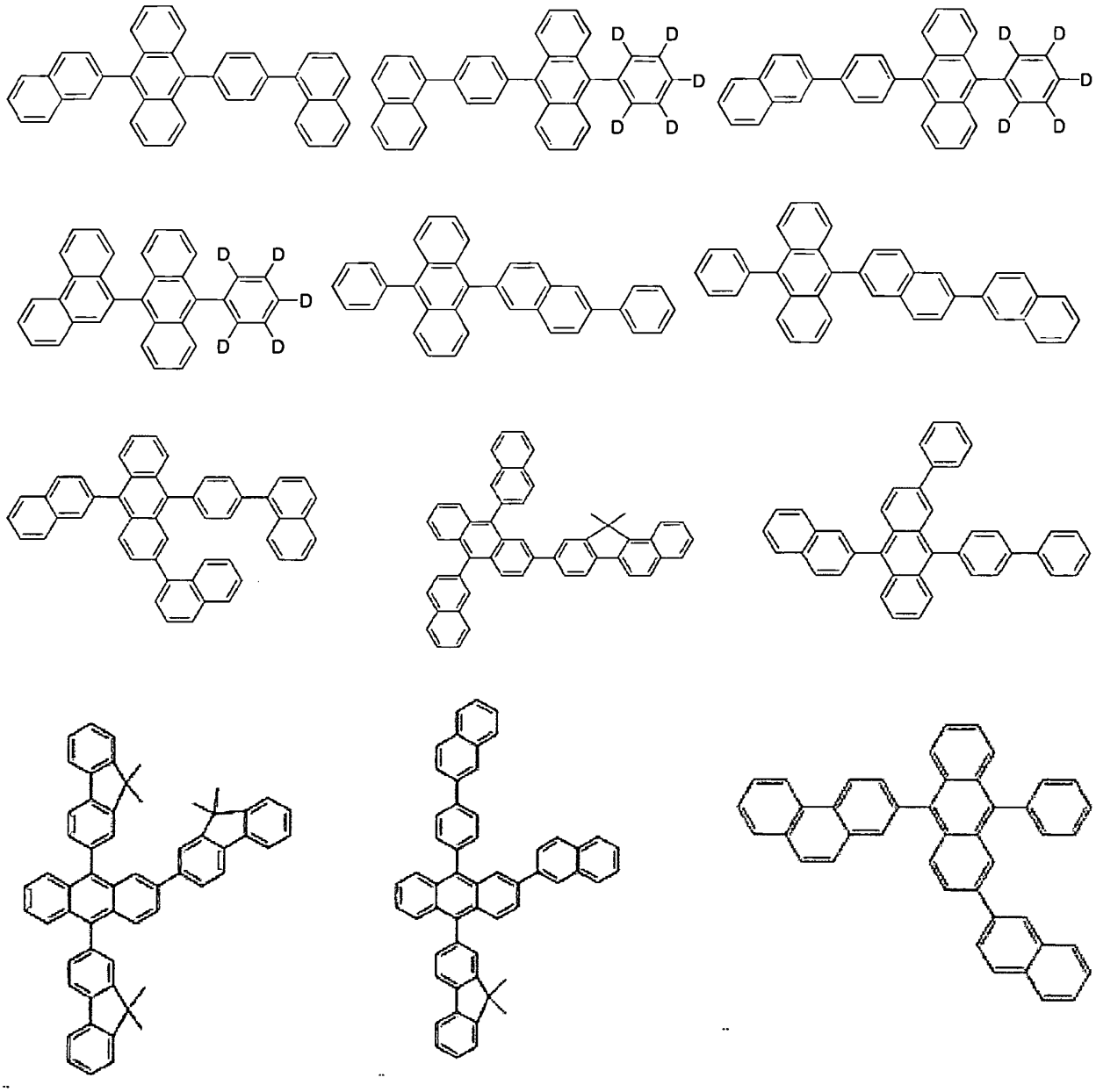






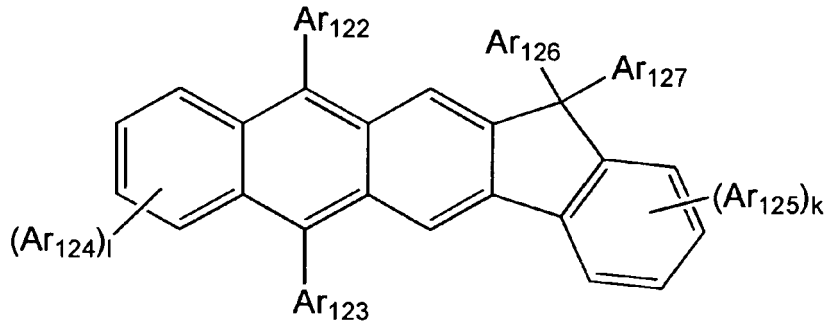






[0106] An anthracene-based compound represented by Formula 401 below may be used as the host.

Formula 401

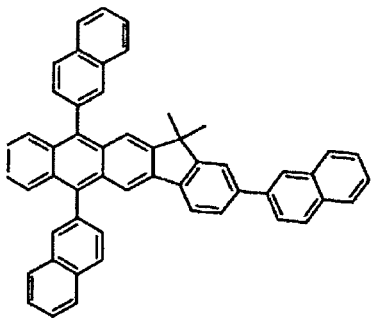


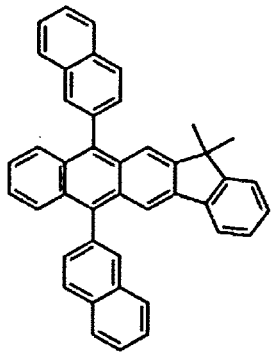
[0107] Ar_{122} to Ar_{125} in Formula 401 above may be defined as described above in conjunction with Ar_{113} of Formula 400, and thus a detailed description thereof will not be provided here.

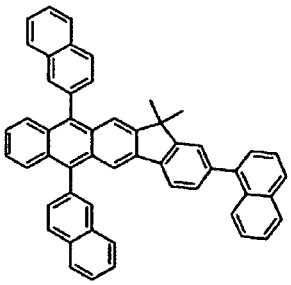
[0108] Ar_{126} and Ar_{127} in Formula 401 above may each independently be a C_1-C_{10} alkyl group, for example, a methyl group, an ethyl group, or a propyl group.

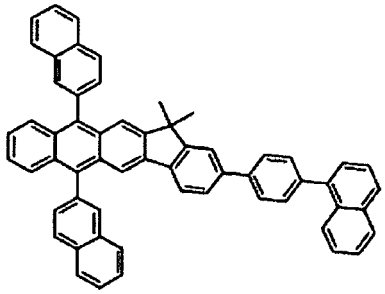
[0109] In Formula 401, k and 1 may each independently be an integer from o to 4, for example, o, 1, or 2.

[0110] For example, the anthracene compound of Formula 401 above may be one of the compounds represented by the following formulae, but is not limited thereto:



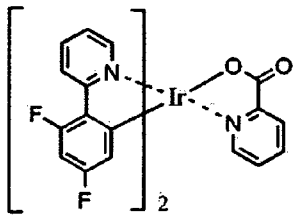




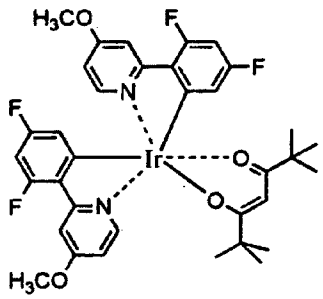


[0111] The amine-based compound in the EML may serve as a host. The EML may further include a dopant, for example, a blue dopant, a green dopant, or a red dopant, in addition to the amine-based compound.

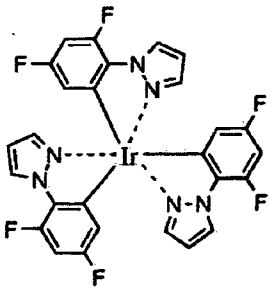
[0112] Non-limiting examples of the blue dopant are compounds represented by the following formulae.



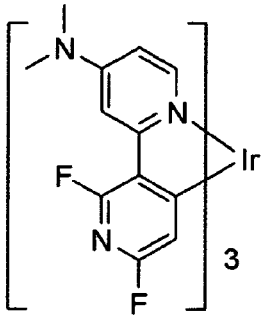
F_2Irpic

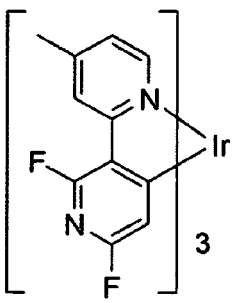


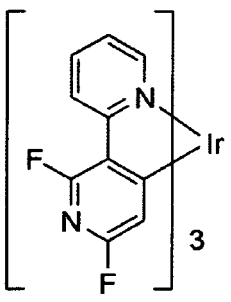
(F2ppy)2Ir(tmd)

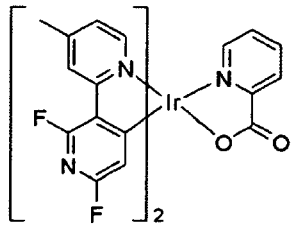


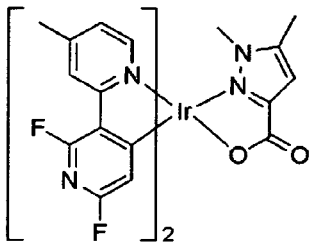
Ir(dfppz)$_3$

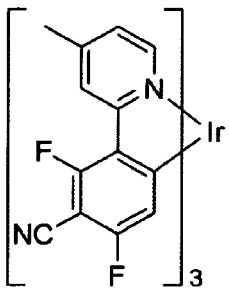


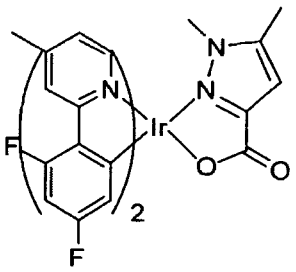


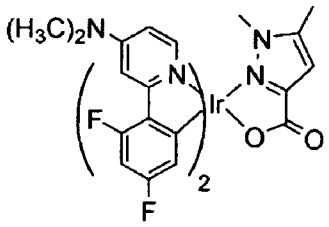


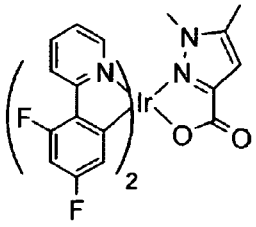


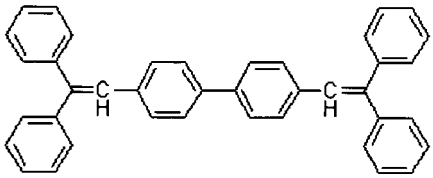




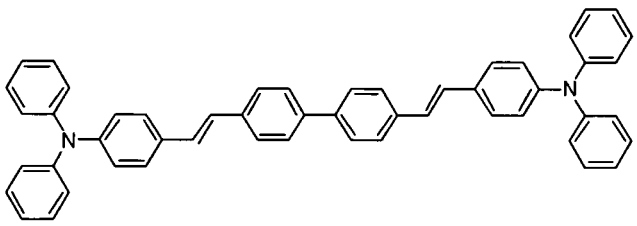




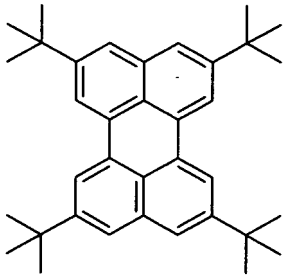




DPVBi

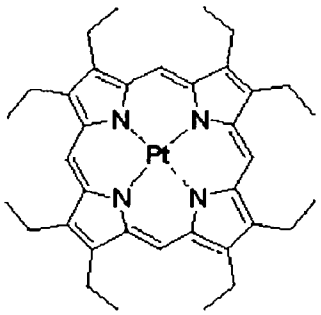


DPAVBi

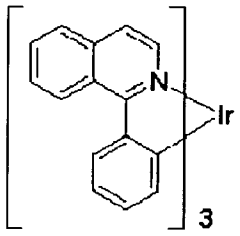


TBPe

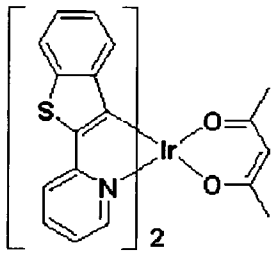
[0113] Non-limiting examples of the red dopant are compounds represented by the following formulae. The red dopant may be DCM or DCJTB, which will be described later.



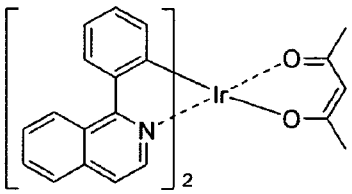
PtOEP

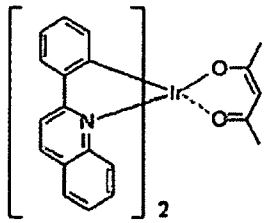


$Ir(piq)_3$

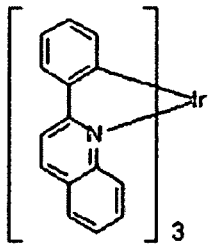


$Btp_2Ir(acac)$

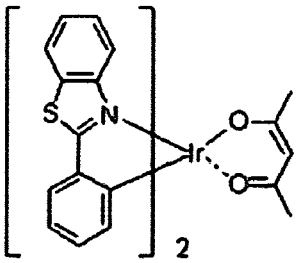




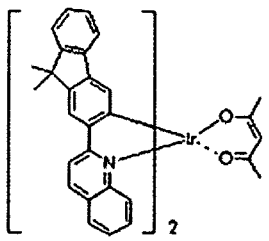
$Ir(pq)_2(acac)$



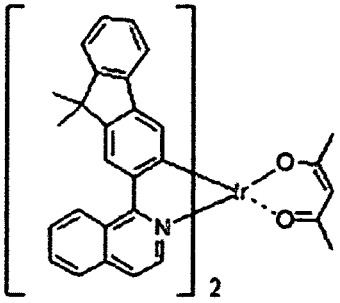
$Ir(2\text{-}phq)_3$



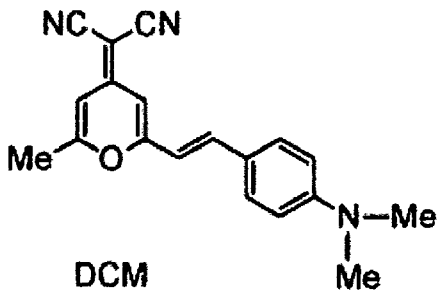
$Ir(BT)_2(acac)$



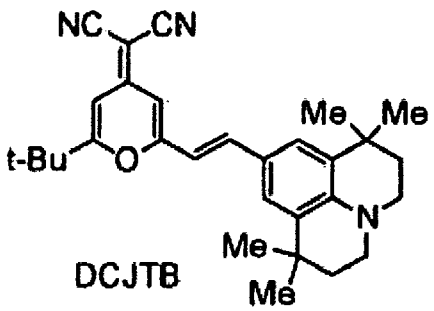
$Ir(flq)_2(acac)$



$Ir(fliq)_2(acac)$

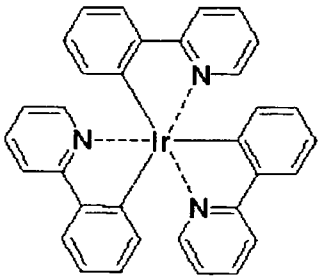


DCM

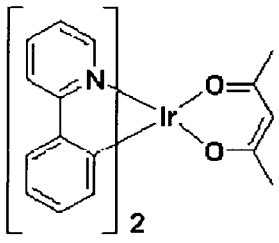


DCJTB

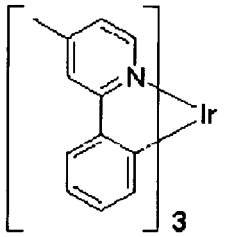
[0114] Non-limiting examples of the green dopant are compounds represented by the following formulae. In an embodiment, the green dopant may be $C_{545}T$ represented below.



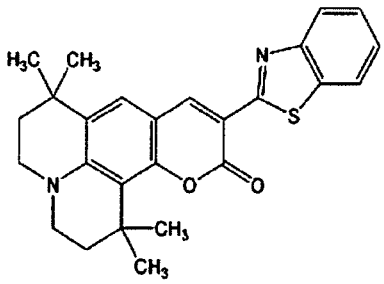
$Ir(ppy)_3$



$Ir(ppy)_2(acac)$

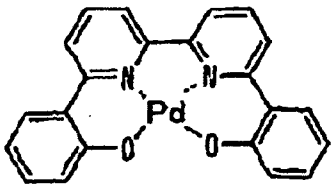


$Ir(mpyp)_3$

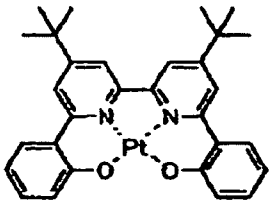


C545T

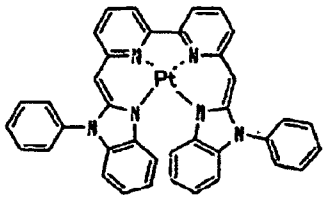
[0115] Non-limiting examples of the dopant that may be used in the EML are Pt complexes represented by the following formulae.



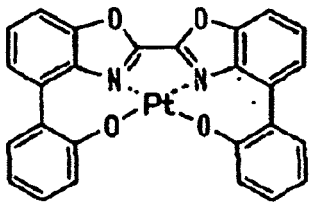
D1



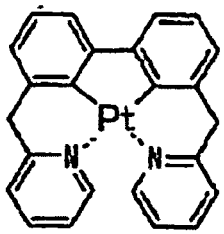
D2



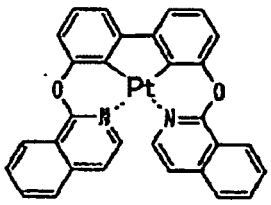
D3



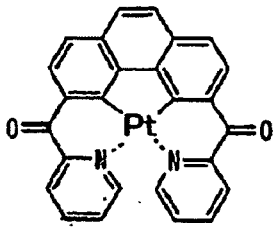
D4



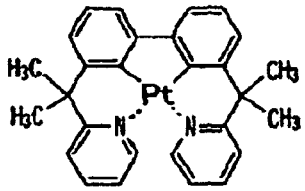
D5



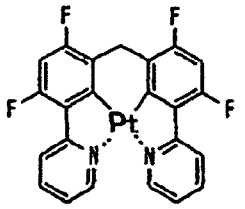
D6



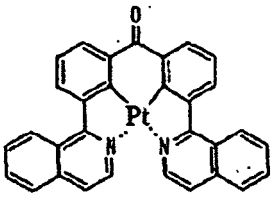
D7



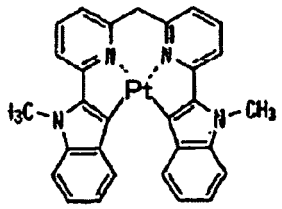
D8



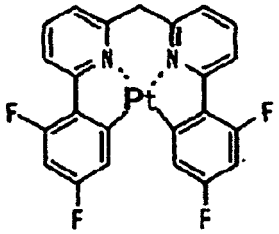
D9



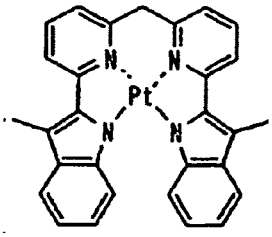
D10



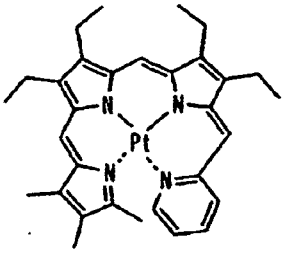
D11



D12



D13



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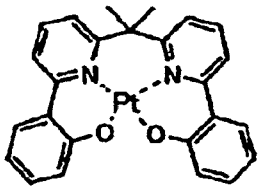
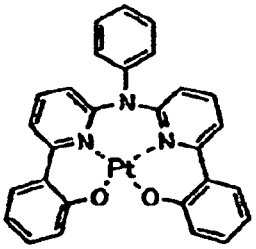
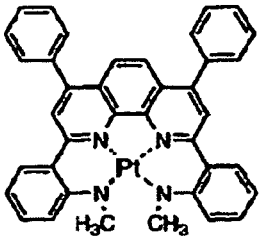


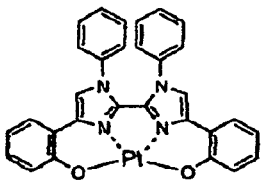
D15



D16



D17



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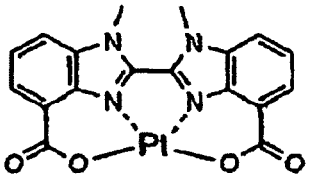

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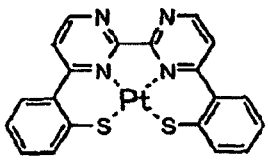
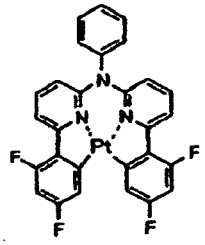
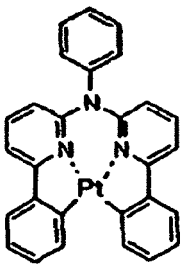


D20



D21



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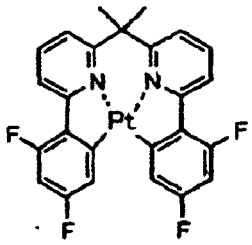

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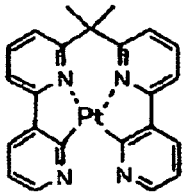

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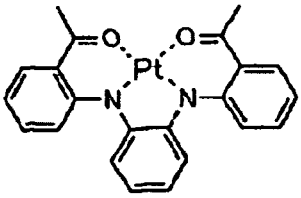

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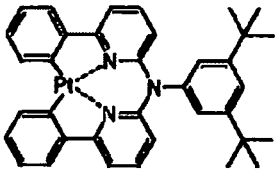

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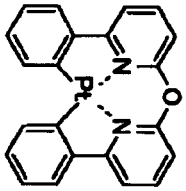

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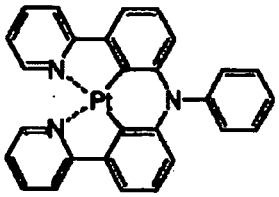

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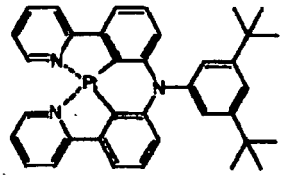

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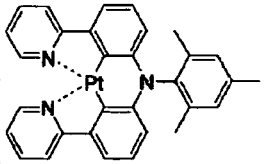

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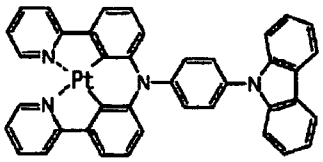
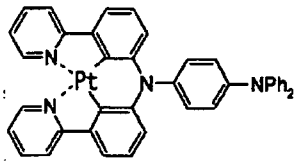


D31



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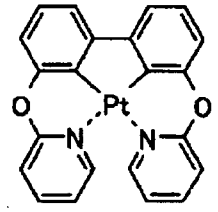

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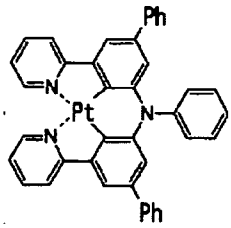

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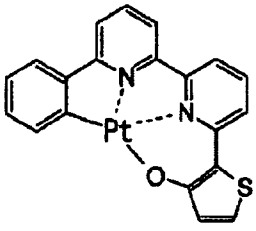
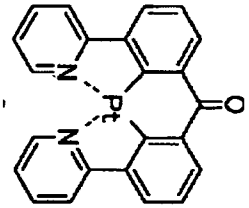
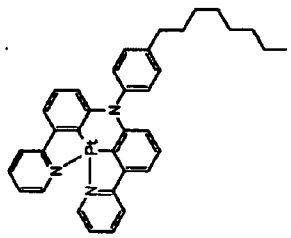


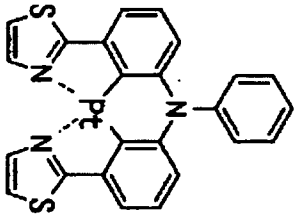
D35



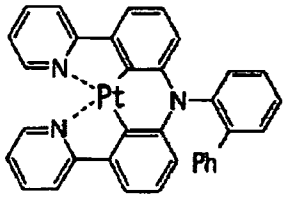
D36



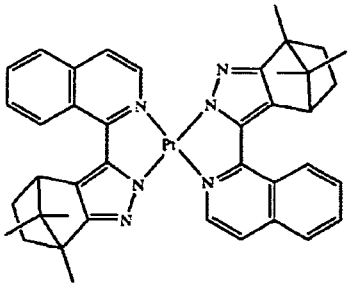
D37



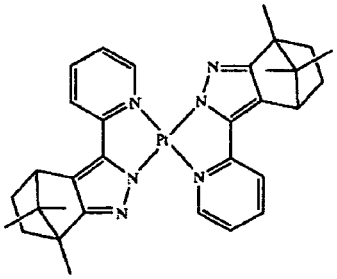
D38



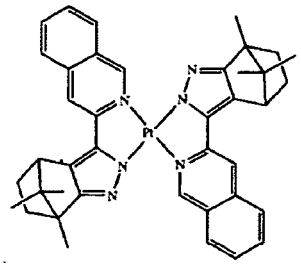
D39



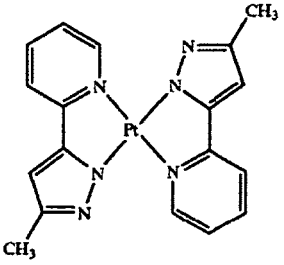
D40



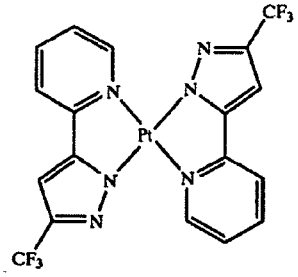
D41



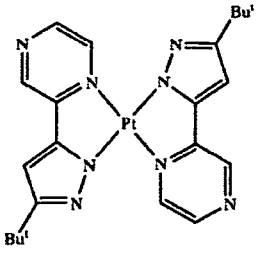
D42



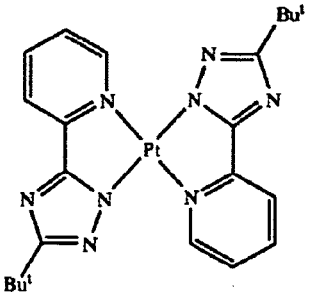
D43



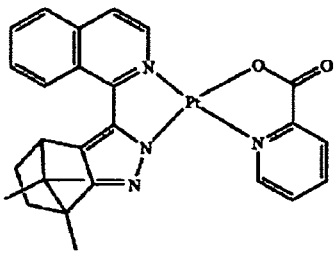
D44



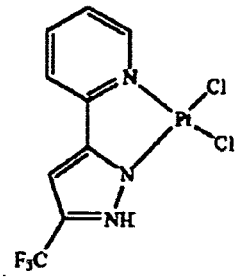
D45



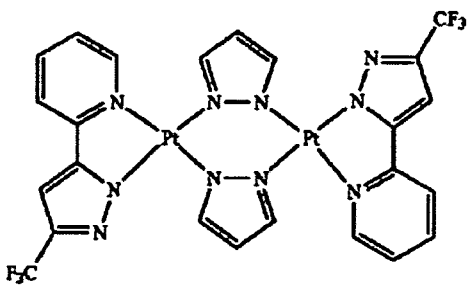
D46



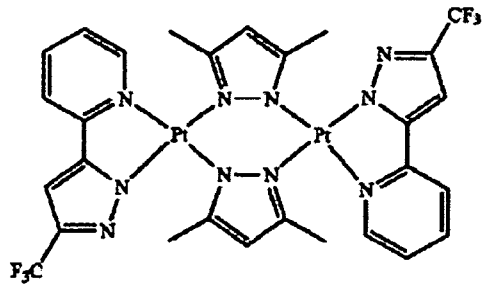
D47



D48

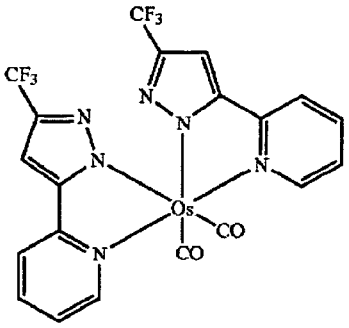


D49

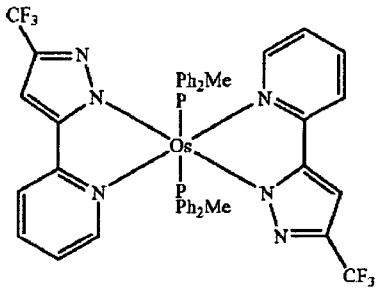


D50

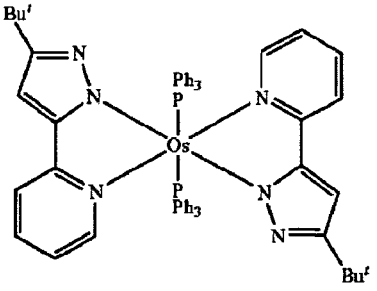
[0116] Non-limiting examples of the dopant that may be used in the EML are Os complexes represented by the following formulae.



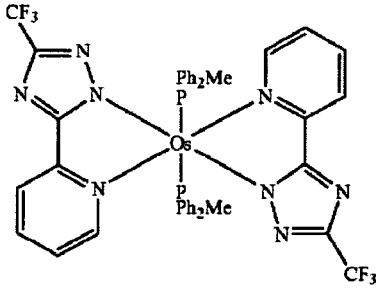
$Os(fppz)_2(CO)_2$



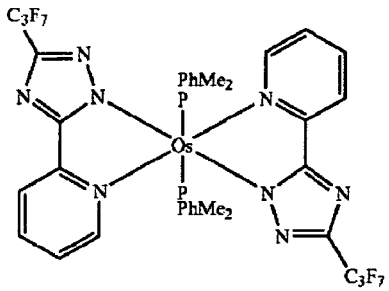
$Os(fppz)_2(PPh_2Me)_2$



$Os(bppz)_2(PPh_3)_2$



$Os(fptz)_2(PPh_2Me)_2$

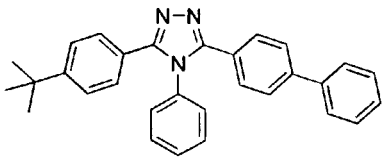


$Os(hptz)_2(PPhMe_2)_2$

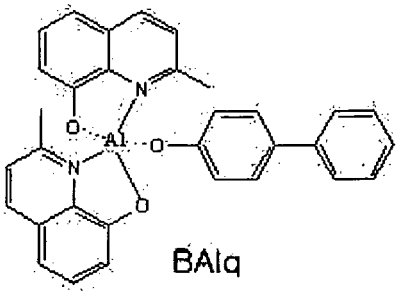
[0117] When the EML includes both a host and a dopant, the amount of the dopant may be from about 0.01 to about 15 parts by weight based on 100 parts by weight of the host. However, the amount of the dopant is not limited to this range.

[0118] The thickness of the EML may be about 10 nm (100 Å) to about 100 nm (1000 Å), and preferably may be about 20 nm (200 Å) to about 60 nm (600 Å). When the thickness of the EML is within these ranges, the EML may have good light emitting ability without a substantial increase in driving voltage.

[0119] Then, an ETL may be formed on the EML by vacuum deposition, spin coating, casting, or the like. When the ETL is formed using vacuum deposition or spin coating, the deposition and coating conditions may be similar to those for the formation of the HIL, though the deposition and coating conditions may vary according to a compound that is used to form the ETL. A material for forming the ETL may be any known material that can stably transport electrons injected from an electron injecting electrode (cathode). Examples of materials for forming the ETL are a quinoline derivative, such as tris(8-quinolinorate)aluminium (Alq_3), TAZ, BAlq, beryllium bis(benzoquinolin-10-olate ($Bebq_2$), 9,10-di(naphthalene-2-yl)anthracene (ADN), Compound 201, and Compound 202, but are not limited thereto.



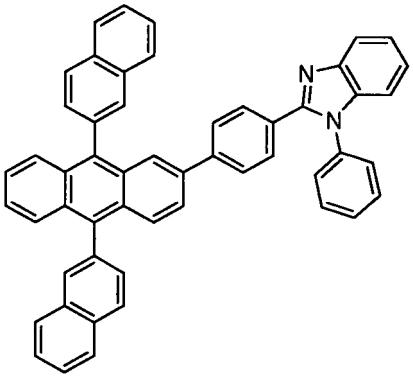
TAZ

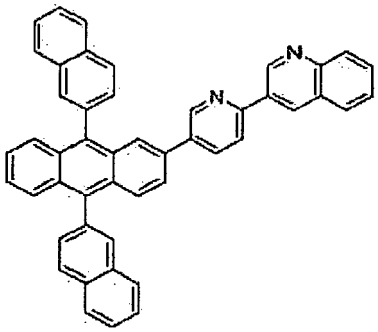


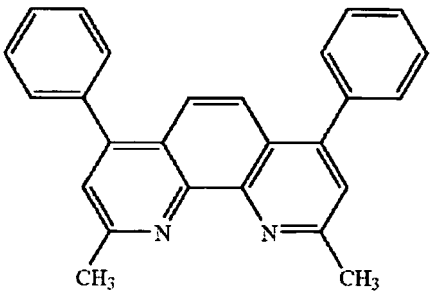
BAlq

Compound 201

Compound 202







BCP

[0120] The ETL may include at least one of the amine-based compounds described above.

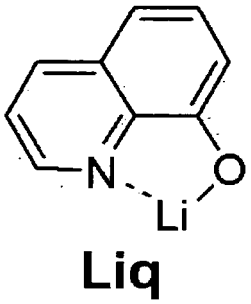
[0121] When the amine-based material of Formula 1 is used as a material for forming the ETL, efficiency and/or lifetime of the organic light-emitting diode may be improved. The ETL including the amine-based compound of Formula 1 may further include a metal complex, for example, lithium quinolate.

[0122] The thickness of the ETL may be in a range of about 10 nm (100 Å) to about 100 nm (1,000 Å), and preferably, may be about 15 nm (150 Å) to about 50 nm (500 Å). When the thickness of the ETL is within these ranges, the ETL

may have satisfactory electron transporting ability without a substantial increase in driving voltage.

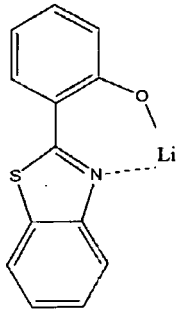
[0123] The ETL may further include a metal-containing material, in addition to any known electron-transporting organic compound.

[0124] The metal-containing material may be a lithium (Li) complex. Non-limiting examples of the Li complex are lithium quinolate (Liq) and Compound 203 below:



Liq

Compound 203



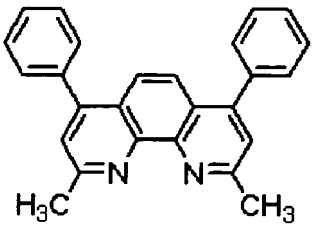
[0125] Then, an EIL, which facilitates injection of electrons from the cathode, may be formed on the ETL. Any suitable electron-injecting material may be used to form the EIL.

[0126] Examples of materials for forming the EIL are LiF, NaCl, CsF, Li_2O, and BaO, which are known in the art. The deposition and coating conditions for forming the EIL 18 may be similar to those for the formation of the HIL, though the deposition and coating conditions may vary according to the material that is used to form the EIL 18.

[0127] The thickness of the EIL may be about 0.1 nm (1 Å) to about 10 nm (100 Å), and preferably, may be about 0.3 nm (3 Å) to about 9 nm (90 Å). When the thickness of the EIL is within these ranges, the EIL may have satisfactory electron injection ability without a substantial increase in driving voltage.

[0128] The second electrode 17 is disposed on the organic layer 15. The second electrode 17 may be a cathode that is an electron injection electrode. A material for forming the second electrode 17 may be a metal, an alloy, an electro-conductive compound, which have a low work function, or a mixture thereof. In this regard, the second electrode 17 may comprise lithium (Li), magnesium (Mg), aluminium (Al), aluminium (Al)-lithium (Li), calcium (Ca), magnesium (Mg)-indium (In), magnesium (Mg)-silver (Ag), or the like, and may be formed as a thin film type transmission electrode. To manufacture a top-emission light-emitting diode, the transmission electrode may comprise indium tin oxide (ITO) or indium zinc oxide (IZO).

[0129] When a phosphorescent dopant is used in the EML, a HBL may be formed between the HTL and the EML or between the H-functional layer and the EML by using vacuum deposition, spin coating, casting, Langmuir-Blodgett (LB) deposition, or the like, in order to prevent diffusion of triplet excitons or holes into the ETL. When the HBL is formed using vacuum deposition or spin coating, the conditions for deposition and coating may be similar to those for the formation of the HIL, although the conditions for deposition and coating may vary according to the material that is used to form the HBL. Any known hole-blocking material may be used. Non-limiting examples of hole-blocking materials are oxadiazole derivatives, triazole derivatives, and phenanthroline derivatives. For example, BCP represented by the following formula may be used as a material for forming the HBL.



BCP

[0130] The thickness of the HBL may be from about 2 nm (20 Å) to about 100 nm (1000 Å), and preferably, may be

from about 3 nm (30 Å) to about 30 nm (300 Å). When the thickness of the HBL is within these ranges, the HBL may have a good hole blocking ability without a substantial increase in driving voltage.

[0131] Examples of the unsubstituted C_1-C_{60} alkyl group (or C_1-C_{60} alkyl group) as used herein are C_1-C_{60} linear or branched alkyl groups, such as methyl, ethyl group, propyl, ethyl, propyl, isobutyl, sec-butyl, pentyl, iso-amyl, and hexyl. Examples of the substituted C_1-C_{60} alkyl group are a C_1-C_{60} alkyl group of which at least one hydrogen atom is substituted with at least one of a deuterium atom; -F; -Cl; -Br; -I; -CN; a hydroxyl group; $-NO_2$; an amino group; an amidino group; hydrazine; hydrazone; a carboxyl group or a salt thereof; a sulfonic acid group or a salt thereof; a phosphoric acid or a salt thereof; a tri(C_6-C_{60}aryl)silyl group; a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, and a C_2-C_{60} alkynyl group; an a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, and a C_2-C_{60} alkynyl group of which at least one hydrogen atom are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, and a phosphoric acid or a salt thereof,; a C_3-C_{60} cycloalkyl group, a C_3-C_{60} cycloalkenyl group, a C_6-C_{60} aryl group, a C_2-C_{60} heteroaryl group, a C_6-C_{60} aralkyl group, a C_6-C_{60} aryloxy group, and a C_6-C_{60} arylthiol group; and a C_3-C_{60} cycloalkyl group, a C_3-C_{60} cycloalkenyl group, a C_6-C_{60} aryl group, a C_2-C_{60} heteroaryl group, a C_6-C_{60} aralkyl group, a C_6-C_{60} aryloxy group, and a C_6-C_{60} arylthiol group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one -F, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group, and a C_2-C_{60} heteroaryl group.

[0132] The unsubstituted C_1-C_{60} alkoxy group (or C_1-C_{60} alkoxy group) may be a group represented by -OA, wherein A is an unsubstituted C_1-C_{60} alkyl group described above. Examples of the unsubstituted C_1-C_{60} alkoxy group are a methoxy group, an ethoxy group, and an isopropyloxy group. At least one of the hydrogen atoms in the alkoxy group may be substituted with the substituents described above in conjunction with the substituted C_1-C_{60} alkyl group.

[0133] The unsubstituted C_2-C_{60} alkenyl group (or C_2-C_{60} alkenyl group) is a hydrocarbon chain having a carbon-carbon double bond in the center or at a terminal of the unsubstituted C_2-C_{60} alkyl group. Examples of the alkenyl group are an ethenyl group, a propenyl group, a butenyl group, and the like. At least one hydrogen atom in the unsubstituted C_2-C_{60} alkenyl group may be substituted with those substituents described in conjunction with the substituted C_1-C_{60} alkyl group.

[0134] The unsubstituted C_2-C_{60} alkynyl group (C_2-C_{60} alkynyl group) is a C_2-C_{60} alkyl group having at least one carbon-carbon triple bond in the center or at a terminal thereof. Examples of the unsubstituted C_2-C_{60} alkynyl group (or C_2-C_{60} alkynyl group) are an ethenyl group, a propynyl group, and the like. At least one hydrogen atom in the alkynyl group may be substituted with the substituents described above in conjunction with the C_1-C_{60} alkyl group.

[0135] As used herein, the unsubstituted C_3-C_{60} cycloalkyl group indicates a cyclic, monovalent C_3-C_{60} saturated hydrocarbon group. Non-limiting examples of the unsubstituted C_3-C_{60} cycloalkyl group are cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl, and cyclooctyl. At least one hydrogen atom in the cyclo alkynyl group may be substituted with those substituents described above in conjunction with the C_1-C_{60} alkyl group.

[0136] As used herein, the unsubstituted C_3-C_{60} cycloalkenyl group indicates a non-aromatic, cyclic unsaturated hydrocarbon group with at least one carbon-carbon double bond. Examples of the unsubstituted C_3-C_{60} cycloalkenyl group are cyclopropenyl, cyclobutenyl, cyclopentenyl, cyclohexenyl, cycloheptenyl, 1,3-cyclohexadienyl, 1,4-cyclohexadienyl, 2,4-cycloheptadienyl, and 1,5-cyclooctadienyl. At least one hydrogen atom in the cycloalkynyl group may be substituted with those substituents described above in conjunction with the substituted C_1-C_{60} alkyl group.

[0137] The unsubstituted C_6-C_{60} aryl group is a monovalent group having a carbocyclic system having 6 to 60 carbon atoms including at least one aromatic ring. The unsubstituted C_6-C_{60} arylene group is a bivalent group having a carbocyclic system having 6 to 60 carbon atoms including at least one aromatic ring. When the aryl group and the arylene group have at least two rings, they may be fused to each other via a single bond. At least one hydrogen atom in the aryl group and the arylene group may be substituted with the substituents described above in conjunction with the C_1-C_{60} alkyl group.

[0138] Examples of the substituted or unsubstituted C_6-C_{60} aryl group are a phenyl group, a C_1-C_{10} alkylphenyl group (e.g., an ethylphenyl group), a C_1-C_{10} alkylbiphenyl group (e.g., an ethylbiphenyl group), a halophenyl group (e.g., an o-, m- or p-fluorophenyl group and a dichlorophenyl group), a dicyanophenyl group, a trifluoromethoxyphenyl group, an o-, m- or p-tolyl group, an o-, m- or p-cumenyl group, a mesityl group, a phenoxyphenyl group, a (α,α-dimethylbenzene)phenyl group, a (N,N'-dimethyl)aminophenyl group, a (N,N'-diphenyl)aminophenyl group, a pentalenyl group, an indenyl group, a naphthyl group, a halonaphthyl group (e.g., a fluoronaphthyl group), a C_1-C_{10} alkylnaphthyl group (e.g., a methylnaphthyl group), a C_1-C_{10} alkoxynaphthyl group (e.g., a methoxynaphthyl group), an anthracenyl group, an azulenyl group, a heptalenyl group, an acenaphthylenyl group, a phenalenyl group, a fluorenyl group, an anthraquinolyl group, a methylanthryl group, a phenanthryl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, an ethylchrysenyl group, a picenyl group, a perylenyl group, a chloroperylenyl group, a pentaphenyl group, a pentacenyl group, a tetraphenylenyl group, a hexaphenyl group, hexacenyl group, a rubicenyl group, a coronenyl group, a trinaphthylenyl group, a heptaphenyl group, a heptacenyl group, a pyranthrenyl group, and an ovalenyl group. Examples of the substituted

C_6-C_{60} aryl group may be inferred based on those of the unsubstituted C_6-C_{60} aryl group and the substituted C_1-C_{60} alkyl group described above. Examples of the substituted or unsubstituted C_6-C_{60} arylene group may be inferred based on those examples of the substituted or unsubstituted C_6-C_{60} aryl group described above.

[0139] The unsubstituted C_2-C_{60} heteroaryl group is a monovalent group having at least one aromatic ring having at least one heteroatom selected from among N, O, P, and S as a ring-forming atom. The unsubstituted C_2-C_{60} heteroarylene group is a bivalent group having at least one aromatic ring having at least one heteroatom selected from among N, O, P, and S. In this regard, when the heteroaryl group and the heteroarylene group have at least two rings, they may be fused to each other via a single bond. At least one hydrogen atom in the heteroaryl group and the heteroarylene group may be substituted with those substituents described above in conjunction with the C_1-C_{60} alkyl group.

[0140] Examples of the unsubstituted C_2-C_{60} heteroaryl group are a pyrazolyl group, an imidazolyl group, an oxazolyl group, a thiazolyl group, a triazolyl group, a tetrazolyl group, an oxadiazolyl group, a pyridinyl group, a pyridazinyl group, a pyrimidinyl group, a triazinyl group, a carbazolyl group, an indolyl group, a quinolinyl group, an isoquinolinyl group, a benzoimidazolyl group, an imidazopyridinyl group, an imidazopyrimidinyl group, a dibenzothiophenyl group, a dibenzofuranyl group and a phenanthrolinyl group. Examples of the unsubstituted C_2-C_{60} heteroarylene group may be inferred based on those examples of the substituted or unsubstituted C_2-C_{60} arylene group described above.

[0141] The substituted or unsubstituted C_6-C_{60} aryloxy group indicates $-OA_2$ (wherein A_2 is a substituted or unsubstituted C_6-C_{60} aryl group described above). The substituted or unsubstituted C_6-C_{60} arylthiol group indicates $-SA_3$ (wherein A_3 is a substituted or unsubstituted C_6-C_{60} aryl group described above).

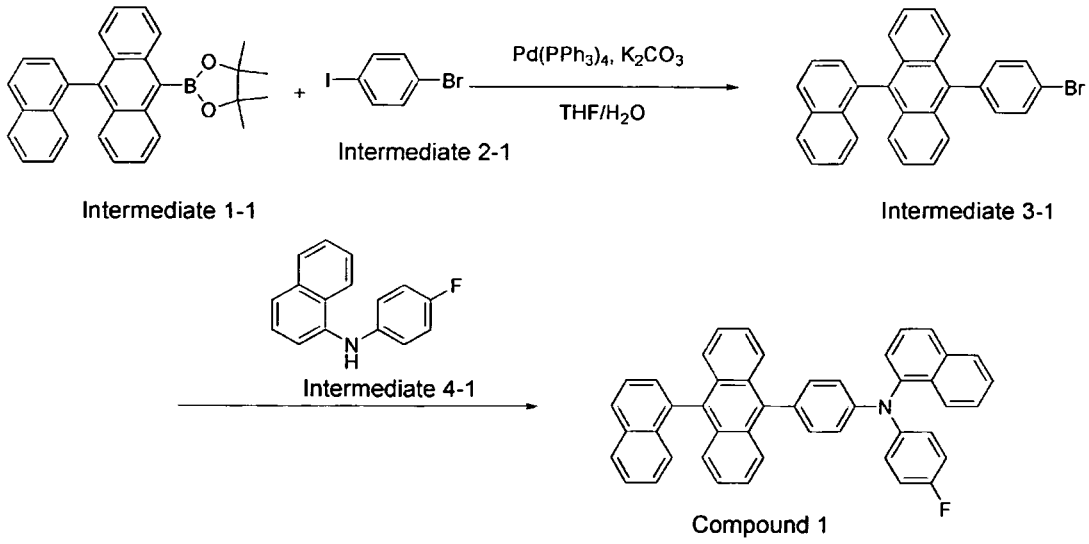
[0142] Hereinafter, the present invention will be described in detail with reference to the following synthesis examples and other examples. However, these examples are for illustrative purposes only and are not intended to limit the scope of the present invention.

EXAMPLES

Synthesis Example 1: Synthesis of Compound 1

[0143] Compound 1 was synthesized according to Reaction Scheme 1 below:

Reaction Scheme 1



Synthesis of Intermediate 3-1

[0144] 8.60g (20.0 mmol) of Intermediate 1-1, 5.66g (20.0 mmol) of Intermediate 2-1, 1.15 g (1.0 mmol) of tetrakis(triphenylphosphine)palladium(o) ($Pd(PPh_3)_4$), and 8.29 g (60.0 mmol) of K_2CO_3 were dissolved in 50 mL of a mixed solution of THF/H_2O (2:1), and then the resultant solution was stirred at about 70°C for about 5 hours. The resultant mixture was cooled to room temperature, followed by three times of extraction with 50 mL of water and 50 mL of diethylether. An organic layer was collected and was dried using magnesium sulfate to evaporate the solvent. The residue was separated and purified by silica gel column chromatography to obtain 7.33 g (80% Yield) of Intermediate 3-1.

Synthesis of Compound 1

[0145] 4.58 g (10.0 mmol) of Intermediate 3-1, 2.85 g (12.0 mmol) of Intermediate 4-1, 0.18 g (0.2 mmol) of $Pd_2(dba)_3$ (tris(dibenzylidineacetone)dipalladium(o)), 0.04 g (0.4 mmol) of tri-tert-butylphosphine ($P(t\text{-}Bu)_3$), and 1.44 g (15.0 mmol) of NaO*t*Bu were dissolved in 50 mL of toluene, and the resultant solution was then refluxed for about 3 hours. The resultant mixture was cooled to room temperature, followed by three times of extraction with 40 mL of water and 40 mL of diethylether. An organic layer was collected and was dried using magnesium sulfate to evaporate the solvent. The residue was separated and purified by silica gel column chromatography to obtain 4.80 g (78% Yield) of Compound 1. This compound was identified using mass spectroscopy/fast atom bombardment (MS/FAB) and ^{1}H NMR.

$C_{46}H_{30}FN$: calc. 615.24, and found; 615.22

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.13-8.11 (dd, 1H), 7.87-7.85 (m, 1H), 7.84-7.80 (m, 3H), 7.72-7.69 (m, 2H), 7.67 (d, 1H), 7.65 (d, 1H), 7.59-7.56 (m, 2H), 7.54-7.51 (dd, 1H), 7.48-7.41 (m, 4H), 7.37-7.23 (m, 6H), 7.18-7.14 (m, 2H), 7.09-7.06 (m, 1H), 6.98-6.94 (m, 2H), 6.85-6.83 (dd, 1H), 6.79-6.75 (m, 2H)

Synthesis Example 2: Synthesis of Compound 3

[0146] 4.98 g of Compound 3 (Yield 80%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-3 instead of Intermediate 4-1 was used. Compound 3 was identified using MS/FAB and ^{1}H NMR.

$C_{47}H_{30}N_2$: calc. 622.24, and found; 622.23

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.84-7.82 (m, 2H), 7.81 (d, 1H), 7.78-7.76 (m, 1H), 7.72-7.68 (m, 3H), 7.66 (d, 1H), 7.65 (d, 1H), 7.63-7.59 (m, 2H), 7.57-7.52 (m, 3H), 7.47-7.43 (m, 3H), 7.41-7.38 (m, 2H), 7.37-7.27 (m, 6H), 7.17 (dd, 1H), 7.13-7.09 (m, 2H), 6.99-6.95 (m, 1H), 6.88-6.85 (m, 1H)

Synthesis Example 3: Synthesis of Compound 4

[0147] 4.86 g of Compound 4 (Yield 75%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-4 instead of Intermediate 4-1 was used. Compound 4 was identified using MS/FAB and ^{1}H NMR.

$C_{49}H_{32}N_2$: calc. 648.26, and found; 648.27

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.86-7.82 (m, 2H), 7.81 (d, 1H), 7.73-7.68 (m, 2H), 7.66 (d, 1H), 7.65-7.52 (m, 3H), 7.60-7.58 (m, 2H) 7.54-7.49 (m, 3H), 7.46-7.42 (m, 3H), 7.40-7.28 (m, 8H), 6.99-6.95 (m, 1H), 6.90-6.84 (m, 4H), 6.74-6.70 (m, 2H)

Synthesis Example 4: Synthesis of Compound 5

[0148] 4.77g of Compound 5 (Yield 70%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-5 instead of Intermediate 4-1 was used. Compound 5 was identified using MS/FAB and ^{1}H NMR.

$C_{51}H_{36}FN$: calc. 681.28, and found; 681.27

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.84-7.82 (m, 2H), 7.80 (d, 1H), 7.78-7.75 (m, 1H), 7.72-7.68 (m, 2H), 7.67 (d, 1H), 7.65 (d, 1H), 7.62-7.58 (m, 2H), 7.56-7.52 (m, 2H), 7.47-7.44 (m, 1H), 7.38-7.27 (m, 6H), 7.14-7.08 (m, 2H), 6.98-6.96 (m, 1H), 6.94-6.89 (m, 2H), 6.85-6.83 (dd, 1H), 6.79-6.77 (m, 2H), 6.75 (d, 1H), 6.73-6.70 (m, 2H), 1.66 (s, 6H)

Synthesis Example 5: Synthesis of Compound 6

[0149] 4.87 g of Compound 6 (Yield 66%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-6 instead of Intermediate 4-1 was used. Compound 6 was identified using MS/FAB and ^{1}H NMR.

$C_{55}H_{35}N_3$: calc. 737.28, and found; 737.29

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.07-8.05 (m, 1H), 7.86-7.83 (m, 2H), 7.81 (d, 1H), 7.73-7.69 (m, 2H), 7.68-7.64 (m, 2H), 7.62-7.58 (m, 2H), 7.50-7.44 (m, 6H), 7.42-7.28 (m, 11H), 7.26-7.23 (m, 2H), 6.99-6.96 (m, 1H), 6.90-6.86 (m, 2H), 6.81-6.78 (dd, 1H), 6.73-6.69 (m, 2H)

Synthesis Example 6: Synthesis of Compound 7

[0150] 4.82 g of Compound 7 (Yield 71%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-7 instead of Intermediate 4-1 was used. Compound 7 was identified

using MS/FAB and ^{1}H NMR.
$C_{49}H_{30}N_2S$: calc. 678.21; and found 678.22
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.15-8.13 (m, 1H), 8.09-8.06 (m, 1H), 7.84-7.80 (m, 4H), 7.72-7.69 (m, 2H), 7.68-7.64 (m, 2H), 7.61-7.58 (m, 2H), 7.54-7.51 (m, 1H), 7.47-7.41 (m, 2H), 7.39-7.27 (m, 8H), 7.16 (d, 1H), 7.13-7.10 (dd, 1H), 7.04-7.01 (m, 1H), 6.93-6.90 (m, 2H), 6.88-6.84 (m, 2H)

Synthesis Example 7: Synthesis of Compound 11

[0151] 4.17 g of Compound 11 (Yield 62%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-11 instead of Intermediate 4-1 was used. Compound 11 was identified using MS/FAB and ^{1}H NMR.
$C_{48}H_{29}F_2NO$: calc. 673.22, and found; 673.21
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.87-7.85 (m, 1H), 7.84-7.81 (m, 3H), 7.76-7.70 (m, 3H), 7.68-7.66 (dd, 1H), 7.65-7.64 (m, 1H), 7.62-7.58 (m, 3H), 7.55-7.50 (m, 3H), 7.46-7.40 (m, 2H), 7.37-7.29 (m, 5H), 7.16-7.13 (dd, 1H), 7.10-7.03 (m, 2H), 6.98-6.93 (m, 2H), 6.86-6.82 (m, 2H)

Synthesis Example 8: Synthesis of Compound 13

[0152] 4.52 g of Compound 13 (Yield 79%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-13 instead of Intermediate 4-1 was used. Compound 13 was identified using MS/FAB and ^{1}H NMR.
$C_{43}H_{28}N_2$: calc. 572.23, and found; 572.23
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.84-7.82 (m, 2H), 7.80 (d, 1H), 7.73-7.70 (m, 2H), 7.67(d, 1H), 7.65 (d, 1H), 7.62-7.58 (m, 2H), 7.54-7.52 (dd, 1H), 7.47-7.43 (m, 2H), 7.38-7.29 (m, 7H), 7.22-7.14 (m, 4H), 7.11-7.06 (m, 1H), 6.97-6.95 (m, 1H), 6.89-6.86 (m, 1H), 6.84-6.81 (m, 2H)

Synthesis Example Q: Synthesis of Compound 14

[0153] 4.29 g of Compound 14 (Yield 75%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-14 instead of Intermediate 4-1 was used. Compound 14 was identified using MS/FAB and 1H NMR.
$C_{43}H_{28}N_2$: calc.572.23, and found; 572.24
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.86-7.82 (m, 2H), 7.81 (d, 1H), 7.73-7.68 (m, 2H), 7.66 (d, 1H), 7.65 (d, 1H), 7.61-7.57 (m, 2H), 7.54-7.49 (m, 1H), 7.46-7.42 (m, 1H), 7.40-7.28 (m, 7H), 7.22-7.17 (m, 4H), 7.10-7.06 (m, 1H), 6.98-6.95 (m, 2H), 6.88-6.84 (m, 1H), 6.80-6.76 (m, 2H)

Synthesis Example 10: Synthesis of Compound 17

[0154] 4.49 g of Compound 17 (Yield 73%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-17 instead of Intermediate 4-1 was used. Compound 17 was identified using MS/FAB and ^{1}H NMR.
$C_{43}H_{28}F_3N$: calc. 615.22, and found; 615.23
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.85-7.81 (m, 2H), 7.79 (d, 1H), 7.74-7.69 (m, 2H), 7.67 (d, 1H), 7.66 (d, 1H), 7.62-7.59 (m, 2H), 7.56-7.48 (m, 3H), 7.43-7.41 (m, 1H), 7.34-7.23 (m, 5H), 7.18-7.15 (m, 4H), 7.06-7.03 (m, 1H), 6.97-6.95 (m, 2H), 6.86-6.83 (m, 1H), 6.78-6.74 (m, 2H)

Synthesis Example 11: Synthesis of Compound 18

[0155] 5.81g of Compound 18 (Yield 70%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-18 instead of Intermediate 4-1 was used. Compound 18 was identified using MS/FAB and ^{1}H NMR.
$C_{61}H_{42}N_2Si$: calc. 830.31, and found; 830.30
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.82-7.80 (m, 2H), 7.79 (d, 1H), 7.73-7.68 (m, 2H), 7.66 (d, 1H), 7.65 (d, 1H), 7.60-7.55 (m, 8H), 7.52-7.49 (m, 1H), 7.45-7.42 (m, 1H), 7.37-7.26 (m, 15H), 7.24-7.20 (m, 3H), 7.16-7.14 (m, 1H), 7.06-7.02 (m, 2H), 6.96-6.94 (m, 2H), 6.80-6.76 (m, 2H)

Synthesis Example 12: Synthesis of Compound 19

[0156] 3.70 g of Compound 19 (Yield 58%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-19 instead of Intermediate 4-1 was used. Compound 19 was identified using MS/FAB and ^{1}H NMR.

$C_{42}H_{24}F_5N$: calc. 637.18, and found; 637.19

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.85-7.83 (m, 2H), 7.81 (d, 1H), 7.75-7.71 (m, 2H), 7.68 (d, 1H), 7.66-62 (m, 3H), 7.56-7.54 (dd, 1H), 7.50-7.46 (m, 1H), 7.40-7.31 (m, 5H), 7.25-7.20 (m, 2H), 7.12-7.09 (m, 1H), 7.02-6.99 (m, 1H), 6.92-6.88 (m, 4H)

Synthesis Example 13: Synthesis of Compound 20

[0157] 4.54 g of Compound 20 (Yield 76%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-20 instead of Intermediate 4-1 was used. Compound 20 was identified using MS/FAB and ^{1}H NMR.

$C_{44}H_{27}N_3$: calc. 597.22, and found; 597.23

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.83-7.81 (m, 2H), 7.80 (d, 1H), 7.72-7.69 (m, 2H), 7.67 (d, 1H), 7.66 (d, 1H), 7.62-7.58 (m, 2H), 7.54-7.51 (m, 1H), 7.47-7.43 (m, 1H), 7.40-7.28 (m, 9H), 7.13-7.10 (m, 1H), 7.02-6.99 (m, 4H), 6.89-6.85 (m, 2H)

Synthesis Example 14: Synthesis of Compound 21

[0158] 3.44 g of Compound 21 (Yield 59%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-21 instead of Intermediate 4-1 was used. Compound 21 was identified using MS/FAB and ^{1}H NMR.

$C_{42}H_{27}F_2N$: calc. 583.21, and found; 583.22

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.84-7.82 (m, 2H), 7.81 (d, 1H), 7.74-7.70 (m, 2H), 7.68 (d, 1H), 7.67 (d, 1H), 7.64-7.61 (m, 2H), 7.55-7.52 (dd, 1H), 7.48-7.44 (m, 1H), 7.37-7.30 (m, 5H), 7.23-7.20 (m, 4H), 7.15-7.09 (m, 5H), 7.04-7.00 (m, 2H)

Synthesis Example 15: Synthesis of Compound 22

[0159] 4.03 g of Compound 22 (Yield 56%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-22 instead of Intermediate 4-1 was used. Compound 22 was identified using MS/FAB and ^{1}H NMR.

$C_{51}H_{33}FN_4$: calc. 720.27, and found; 720.28

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.52-8.47 (m, 4H), 7.87-7.82 (m, 5H), 7.76-7.73 (m, 2H), 7.70 (d, 1H), 7.68 (d, 1H), 7.62-7.56 (m, 5H), 7.50-7.46 (m, 1H), 7.44-7.31 (m, 7H), 7.24-7.18 (m, 2H), 7.13-7.09 (m, 2H), 7.00-6.98 (m, 1H), 6.85-6.82 (m, 2H)

Synthesis Example 16: Synthesis of Compound 23

[0160] 5.35 g of Compound 23 (Yield 70%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-23 instead of Intermediate 4-1 was used. Compound 23 was identified using MS/FAB and ^{1}H NMR.

$C_{56}H_{36}N_4$: calc. 764.29, and found; 764.28

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.85-7.82 (m, 3H), 7.81-7.77 (m, 3H), 7.74-7.70 (m, 2H), 7.69-7.67 (m, 1H), 7.66-7.65 (m, 2H), 7.61-7.52 (m, 5H), 7.47-7.30 (m, 12H), 7.27-7.23 (m, 1H), 7.12-7.10 (m, 1H), 7.02-6.98 (m, 4H), 6.93-6.91 (m, 2H)

Synthesis Example 17: Synthesis of Compound 24

[0161] 4.74 g of Compound 24 (Yield 76%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-24 instead of Intermediate 4-1 was used. Compound 24 was identified using MS/FAB and ^{1}H NMR.

$C_{47}H_{32}N_2$: calc. 624.26, and found; 624.25

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.49-8.46 (m, 1H), 7.84-7.78 (m, 5H), 7.76-7.68 (m, 3H), 7.67-7.64 (m, 3H), 7.62-7.59 (m, 2H), 7.55-7.53 (m, 1H), 7.48-7.45 (m, 1H), 7.39-7.28 (m, 6H), 7.23-7.19 (m, 4H), 7.11-7.09 (m, 1H), 7.02-6.98 (m, 3H), 6.90-6.86 (m, 2H)

Synthesis Example 18: Synthesis of Compound 25

[0162] 4.93 g of Compound 25 (Yield 79%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-25 instead of Intermediate 4-1 was used. Compound 25 was identified using MS/FAB and ^{1}H NMR.

$C_{47}H_{32}N_2$: calc. 624.26, and found; 624.25

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.69 (d, 1H), 8.46-8.42 (m, 1H), 7.91-7.97 (m, 1H), 7.86-7.81 (m, 3H), 7.75-7.70 (m, 2H), 7.69-7.66 (m, 2H), 7.63-7.60 (m, 2H), 7.57-7.54 (m, 1H), 7.51-7.47 (m, 2H), 7.40-7.29 (m, 7H), 7.15-7.10 (m, 4H), 6.98-6.91 (m, 3H), 6.87-6.85 (m, 1H), 6.82-6.78 (m, 2H)

Synthesis Example 19: Synthesis of Compound 26

[0163] 4.99 g of Compound 26 (Yield 80%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-26 instead of Intermediate 4-1 was used. Compound 26 was identified using MS/FAB and ^{1}H NMR.

$C_{47}H_{32}N_2$: calc. 624.26, and found; 624.25

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.56-8.53 (m, 2H), 7.85-7.82 (m, 2H), 7.80 (d, 1H), 7.73-7.69 (m, 2H), 7.68-7.66 (m, 1H), 7.65 (d, 1H), 7.61-7.51 (m, 7H), 7.47-7.44 (m, 1H), 7.38-7.26 (m, 5H), 7.18-7.13 (m, 4H), 7.09-7.04 (m, 1H), 6.93-6.88 (m, 3H), 6.84-6.80 (m, 2H)

Synthesis Example 20: Synthesis of Compound 27

[0164] 4.29 g of Compound 27 (Yield 66%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-27 instead of Intermediate 4-1 was used. Compound 27 was identified using MS/FAB and ^{1}H NMR.

$C_{48}H_{31}N_3$: calc. 649.25, and found; 649.26

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.72 (d, 1H), 8.42-8.39 (m, 1H), 7.95-7.92 (m, 1H), 7.84-7.81 (m, 2H), 7.79 (d, 1H), 7.71-7.68 (m, 2H), 7.66 (d, 1H), 7.64 (d, 1H), 7.62-7.57 (m, 2H), 7.55-7.52 (m, 1H), 7.49-7.44 (m, 2H), 7.40-7.26 (m, 9H), 7.11-7.06 (m, 1H), 7.02-7.00 (m, 2H), 6.93-6.89 (m, 2H), 6.79-6.77 (m, 2H)

Synthesis Example 21: Synthesis of Compound 30

[0165] 4.99 g of Compound 30 (Yield 74%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-30 instead of Intermediate 4-1 was used. Compound 30 was identified using MS/FAB and ^{1}H NMR.

$C_{51}H_{34}N_2$: calc. 674.27, and found; 674.26

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.74 (d, 1H), 8.44-8.31 (m, 1H), 8.12-8.08 (dd, 1H), 7.94-7.90 (m, 1H), 7.87-7.80 (m, 4H), 7.75-7.71 (m, 2H), 7.68 (d, 1H), 7.65 (d, 1H), 7.60-7.56 (m, 2H), 7.54-7.52 (m, 1H), 7.49-7.41 (m, 5H), 7.38-7.21 (m, 8H), 7.11-7.07 (m, 2H), 7.03-7.00 (m, 3H), 6.87-6.85 (dd, 1H)

Synthesis Example 22: Synthesis of Compound 31

[0166] 4.98 g of Compound 31 (Yield 80%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-31 instead of Intermediate 4-1 was used. Compound 31 was identified using MS/FAB and ^{1}H NMR.

$C_{47}H_{30}N_2$: calc. 622.24, and found; 622.23

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.05-8.01 (dd, 1H), 7.89-7.86 (m, 1H), 7.84-7.80 (m, 3H), 7.74-7.69 (m, 2H), 7.67 (d, 1H), 7.65 (d, 1H), 7.59-7.56 (m, 2H), 7.53-7.50 (m, 1H), 7.48-7.41 (m, 4H), 7.40-7.22 (m, 8H), 7.13-7.09 (m, 2H), 7.05-7.03 (m, 1H), 6.96-6.93 (dd, 1H), 6.88-6.84 (m, 2H)

Synthesis Example 23: Synthesis of Compound 32

[0167] 5.40 g of Compound 32 (Yield 73%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-32 instead of Intermediate 4-1 was used. Compound 32 was identified using MS/FAB and ^{1}H NMR.

$C_{56}H_{40}N_2$: calc. 740.32, and found; 740.31

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.76 (d, 1H), 8.51-8.47 (dd, 1H), 7.91-7.88 (m, 1H), 7.83-7.79 (m, 3H), 7.78-7.76 (m, 1H), 7.73-7.69 (m, 2H), 7.68 (d, 1H), 7.66 (d, 1H), 7.62-7.58 (m, 2H), 7.56-7.51 (m, 2H), 7.47-7.43 (m, 2H), 7.39-7.26

(m, 8H), 7.15-7.10 (m, 2H), 7.06-7.04 (m, 1H), 6.94-6.90 (m, 3H), 6.87 (d, 1H), 6.83-6.81 (m, 2H), 1.67 (s, 6H)

Synthesis Example 24: Synthesis of Compound 35

[0168] 5.05 g of Compound 35 (Yield 67%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-35 instead of Intermediate 4-1 was used. Compound 35 was identified using MS/FAB and ^{1}H NMR.
$C_{58}H_{38}N_2O$: calc. 778.30, and found; 778.29
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.84-7.80 (m, 4H), 7.78-7.76 (m, 1H), 7.74-7.65 (m, 6H), 7.62-7.59 (m, 3H), 7.55-7.49 (m, 4H), 7.46-7.40 (m, 3H), 7.38-7.27 (m, 5H), 7.19-7.11 (m, 2H), 7.05 (d, 1H), 6.87 (d, 1H), 6.79-6.74 (m, 2H), 1.64 (s, 6H)

Synthesis Example 25: Synthesis of Compound 36

[0169] 5.80 g of Compound 36 (Yield 70%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-36 instead of Intermediate 4-1 was used. Compound 36 was identified using MS/FAB and ^{1}H NMR.
$C_{63}H_{44}N_2$: calc. 828.35, and found; 828.34
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.08-8.04 (m, 2H), 7.85-7.82 (m, 2H), 7.81 (d, 1H), 7.77-7.75 (m, 1H), 7.72-7.68 (m, 2H), 7.68 (d, 1H), 7.66 (d, 1H), 7.62-7.58 (m, 2H), 7.56-7.52 (m, 2H), 7.47-7.43 (m, 1H), 7.36-7.26 (m, 12H), 7.15-7.10 (m, 4H), 7.04-7.02 (m, 1H), 6.99-6.91 (m, 3H), 6.89 (d, 1H), 6.85-6.81 (m, 2H), 1.65 (s, 6H)

Synthesis Example 26: Synthesis of Compound 18

[0170] 5.04 g of Compound 38 (Yield 75%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate 4-38 instead of Intermediate 4-1 was used. Compound 38 was identified using MS/FAB and ^{1}H NMR.
$C_{51}H_{32}N_2$: calc. 672.26, and found; 672.25
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.50-8.47 (m, 1H), 8.16-8.12 (m, 1H), 7.96-7.92 (m, 1H), 7.86-7.82 (m, 2H), 7.80 (d, 1H), 7.73-7.69 (m, 4H), 7.67 (d, 1H), 7.65 (d, 1H), 7.60-7.52 (m, 5H), 7.47-7.41 (m, 2H), 7.39-7.28 (m, 8H), 7.22-7.18 (m, 2H), 7.07-7.03 (m, 1H), 6.97-6.95 (m, 2H)

Synthesis Example 27: Synthesis of Compound 42

Synthesis of Intermediate 3-42

[0171] Intermediate 3-42 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-42 instead of Intermediate 2-1 was used.

Synthesis of Compound 42

[0172] 4.54 g of Compound 42 (Yield 65%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-42 and 4-27, instead of Intermediates 3-1 and 4-1, were used. Compound 42 was identified using MS/FAB and ^{1}H NMR.
$C_{52}H_{33}N_3$: calc. 699.27, and found; 699.28
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.78 (d, 1H), 8.45-8.42 (m, 1H), 8.11-8.08 (m, 1H), 7.94-7.91 (m, 1H), 7.89-7.86 (m, 1H), 7.83-7.80 (m, 2H), 7,78 (d, 1H), 7.72-7.67 (m, 3H), 7.65-7.62 (dd, 1H), 7.58-7.55 (m, 3H), 7.54-7.52 (m, 1H), 7.48-7.43 (m, 2H), 7.40-7.33 (m, 5H), 7.31-7.27 (m, 4H), 7.18-7.15 (dd, 1H), 6.99-6.95 (m, 1H), 6.82-6.79 (m, 2H), 6.77-6.73 (m, 2H)

Synthesis Example 28: Synthesis of Compound 43

[0173] 5.24 g of Compound 43 (Yield 71%) was prepared in the same manner as in the method of preparing Compound 42 of Synthesis Example 27, except that Intermediate 4-43 instead of Intermediate 4-27 was used. Compound 43 was identified using MS/FAB and ^{1}H NMR.
$C_{56}H_{38}N_2$: calc. 738.30, and found; 738.31
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.10-8.08 (m, 1H), 7.90-7.87 (m, 1H), 7.85-7.82 (m, 2H), 7.80 (d, 1H), 7.78-7.75 (m, 1H), 7.71-7.65 (m, 4H), 7.59-7.57 (m, 1H), 7.55-7.51 (m, 4H), 7.46-7.42 (m, 1H), 7.39-7.27 (m, 8H), 7.16-7.10 (m, 3H), 6.97-6.93 (m, 1H), 6.89-6.86 (m, 2H), 6.83-6.81 (dd, 1H), 6.79 (d, 1H), 1.66 (s, 6H)

Synthesis Example 29: Synthesis of Compound 45

Synthesis of Intermediate 3-45

[0174] Intermediate 3-45 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-45 instead of Intermediate 2-1 was used.

Synthesis of Compound 45

[0175] 5.38 g of Compound 45 (Yield 80%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-45 and 4-31, instead of Intermediates 3-1 and 4-1, were used. Compound 45 was identified using MS/FAB and ^{1}H NMR.
$C_{51}H_{32}N_2$: calc. 672.26, and found; 672.27
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.03-8.00 (m, 1H), 7.88-7.85 (m, 1H), 7.84-7.79 (m, 5H), 7.73-7.68 (m, 2H), 7.65-7.63 (m, 1H), 7.60 (d, 1H), 7.55-7.37 (m, 9H), 7.35-7.28 (m, 5H), 7.20-7.15 (m, 2H), 7.10-7.06 (m, 1H), 6.97-6.92 (m, 1H), 6.86-6.85 (dd, 1H), 6.80-6.77 (m, 2H)

Synthesis Example 30: Synthesis of Compound 48

[0176] 5.50 g of Compound 48 (Yield 76%) was prepared in the same manner as in the method of preparing Compound 45 of Synthesis Example 29, except that Intermediate 4-48 instead of Intermediate 4-31 was used. Compound 48 was identified using MS/FAB and ^{1}H NMR.
$C_{55}H_{36}N_2$: calc. 724.29, and found; 724.30
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.56-8.53 (dd, 1H), 8.06-8.03 (m, 1H), 7.85-7.78 (m, 6H), 7.76-7.69 (m, 5H), 7.66-7.60 (m, 3H), 7.55-7.40 (m, 7H), 7.36-7.32 (m, 5H), 7.29-7.25 (m, 1H), 7.21-7.17 (m, 2H), 7.11-7.07 (m, 1H), 6.99-6.94 (m, 1H), 6.87-6.85 (dd, 1H), 6.82-6.79 (m, 2H)

Synthesis Example 31: Synthesis of Compound 51

Synthesis of Intermediate 3-51

[0177] Intermediate 3-51 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-51 instead of Intermediate 2-1 was used.

Synthesis of Compound 51

[0178] 4.29 g of Compound 51 (Yield 66%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-51 and 4-4, instead of Intermediates 3-1 and 4-1, were used. Compound 51 was identified using MS/FAB and ^{1}H NMR.
$C_{48}H_{31}N_3$: calc. 649.25, and found; 649.26
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.50 (d, 1H), 7.93-7.90 (m, 2H), 7.89-7.87 (m, 2H), 7.85 (d, 1H), 7.84-7.81 (m, 1H), 7.74-7.69 (m, 2H), 7.65-7.61 (m, 2H), 7.56-7.49 (m, 5H), 7.46-7.37 (m, 6H), 7.36-7.34 (m, 1H), 7.32-7.27 (m, 2H), 7.14-7.10 (m, 2H), 6.99-6.94 (m, 2H), 6.83-6.79 (m, 2H)

Synthesis Example 32: Synthesis of Compound 54

[0179] 4.23 g of Compound 54 (Yield 65%) was prepared in the same manner as in the method of preparing Compound 51 of Synthesis Example 31, except that Intermediate 4-27 instead of Intermediate 4-4 was used. Compound 54 was identified using MS/FAB and ^{1}H NMR.
$C_{47}H_{30}N_4$: calc. 650.25, and found; 650.24
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.73 (m, 1H), 8.55 (d, 1H), 8.52-8.47 (m, 1H), 7.96-7.88 (m, 5H), 7.86 (d, 1H), 7.84-7.81 (m, 1H), 7.73-7.68 (m, 2H), 7.56-7.53 (m, 1H), 7.49-7.38 (m, 6H), 7.36-7.28 (m, 5H), 7.16-7.13 (m, 2H), 7.00-6.95 (m, 2H), 6.89-6.85 (m, 2H)

Synthesis Example 33: Synthesis of Compound 57

Synthesis of Intermediate 3-57

[0180] Intermediate 3-57 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-57 instead of Intermediate 2-1 was used.

Synthesis of Compound 57

[0181] 4.97 g of Compound 57 (Yield 67%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-57 and 4-32, instead of Intermediates 3-1 and 4-1, were used. Compound 57 was identified using MS/FAB and ^{1}H NMR.
$C_{55}H_{39}N_3$: calc. 741.31, and found; 741.32
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.77 (dd, 1H), 8.52 (d, 1H), 8.23-8.20 (m, 1H), 7.93-7.90 (m, 1H), 7.89 (d, 1H), 7.87 (d, 1H), 7.85-7.82 (m, 1H), 7.78-7.75 (m, 1H), 7.72-7.68 (m, 2H), 7.53-7.44 (m, 6H), 7.42-7.23 (m, 9H), 7.14-7.09 (m, 2H), 6.99-6.95 (m, 1H), 6.87-6.83 (m, 4H), 6.78 (d, 1H), 1.63 (s, 6H)

Synthesis Example 34: Synthesis of Compound 58

Synthesis of Intermediate 3-58

[0182] Intermediate 3-58 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-58 instead of Intermediate 2-1 was used.

Synthesis of Compound 58

[0183] 5.67 g of Compound 58 (Yield 81%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-58 and 4-25, instead of Intermediates 3-1 and 4-1, were used. Compound 58 was identified using MS/FAB and ^{1}H NMR.
$C_{53}H_{36}N_2$: calc. 700.29, and found; 700.30
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.74 (dd, 1H), 8.51-8.48 (m, 1H), 7.94-7.90 (m, 1H), 7.83-7.77 (m, 5H), 7.74-7.65 (m, 6H), 7.55-7.52 (m, 1H), 7.48-7.44 (m, 4H), 7.38-7.25 (m, 7H), 7.19-7.15 (m, 2H), 7.08-7.03 (m, 3H), 6.96-6.92 (m, 2H), 6.86-6.83 (m, 1H), 6.80-6.79 (m, 2H)

Synthesis Example 35: Synthesis of Compound 59

[0184] 5.46 g of Compound 59 (Yield 78%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-24 instead of Intermediate 4-25 was used. Compound 59 was identified using MS/FAB and ^{1}H NMR.
$C_{53}H_{36}N_2$: calc. 700.29, and found; 700.29
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.58-8.54 (m, 1H), 7.86-7.80 (m, 7H), 7.78-7.70 (m, 5H), 7.68-7.65 (m, 3H), 7.54-7.51 (m, 1H), 7.47-7.44 (m, 3H), 7.37-7.26 (m, 6H), 7.17-7.13 (m, 2H), 7.09-7.07 (m, 1H), 7.01-6.97 (m, 2H), 6.98-6.93 (m, 3H), 6.84-6.81 (m, 2H)

Synthesis Example 36: Synthesis of Compound 60

[0185] 5.08 g of Compound 60 (Yield 70%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-27 instead of Intermediate 4-25 was used. Compound 60 was identified using MS/FAB and ^{1}H NMR.
$C_{54}H_{35}N_3$: calc. 725.28, and found; 725.27
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.71-8.69 (m, 1H), 8.44-8.41 (m, 1H), 7.93-7.91 (m, 1H), 7.84-7.78 (m, 5H), 7.74-7.69 (m, 4H), 7.68-7.65 (m, 2H), 7.55-7.52 (m, 1H), 7.48-7.43 (m, 4H), 7.39-7.24 (m, 9H), 7.12-7.11 (m, 1H), 7.03-6.99 (m, 2H), 6.93-6.89 (m, 2H), 6.82-6.80 (m, 2H)

Synthesis Example 37: Synthesis of Compound 61

[0186] 6.04 g of Compound 61 (Yield 74%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-32 instead of Intermediate 4-25 was used. Compound 61 was

identified using MS/FAB and ^{1}H NMR.
$C_{62}H_{44}N_2$: calc. 816.35, and found; 816.34
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.69-8.67 (dd, 1H), 8.41-8.38 (m, 1H), 7.91-7.88 (m, 1H), 7.83-7.76 (m, 6H), 7.74-7.68 (m, 4H), 7.67-7.64 (m, 2H), 7.56-7.51 (m, 2H), 7.47-7.44 (m, 4H), 7.37-7.25 (m, 8H), 7.15-7.09 (m, 2H), 6.99-6.94 (m, 1H), 6.90-6.86 (m, 3H), 6.83-6.79 (m, 2H), 6.77-6.76 (m, 1H), 1.64 (s, 6H)

Synthesis Example 38: Synthesis of Compound 62

[0187] 5.93 g of Compound 62 (Yield 79%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-30 instead of Intermediate 4-25 was used. Compound 62 was identified using MS/FAB and ^{1}H NMR.
$C_{57}H_{38}N_2$: calc. 750.30, and found; 750.29
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.51-8.48 (m, 2H), 8.17-8.15 (dd, 1H), 7.87-7.78 (m, 6H), 7.74-7.65 (m, 6H), 7.55-7.51 (m, 5H), 7.48-7.40 (m, 6H), 7.38-7.28 (m, 5H), 7.23 (t, 1H), 7.14-7.13 (m, 1H), 7.06-7.02 (m, 2H), 6.95-6.93 (dd, 1H), 6.88-6.85 (m, 2H)

Synthesis Example 39: Synthesis of Compound 63

[0188] 5.85 g of Compound 63 (Yield 78%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-63 instead of Intermediate 4-25 was used. Compound 63 was identified using MS/FAB and ^{1}H NMR.
$C_{57}H_{38}N_2$: calc. 750.30, and found; 750.31
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.50-8.47 (m, 2H), 8.15-8.13 (dd, 1H), 7.86-7.77 (m, 6H), 7.73-7.69 (m, 4H), 7.68-7.65 (m, 2H), 7.56-7.52 (m, 5H), 7.49-7.39 (m, 6H), 7.37-7.29 (m, 5H), 7.22 (t, 1H), 7.18-7.16 (m, 1H), 7.11-7.07 (m, 2H), 6.96-6.94 (dd, 1H), 6.89-6.86 (m, 2H)

Synthesis Example 40: Synthesis of Compound 64

[0189] 5.40 g of Compound 64 (Yield 72%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-48 instead of Intermediate 4-25 was used. Compound 64 was identified using MS/FAB and ^{1}H NMR.
$C_{57}H_{38}N_2$: calc. 750.30, and found; 750.31
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.48-8.45 (m, 1H), 8.16-8.14 (dd, 1H), 7.87-7.82 (m, 3H), 7.81-7.68 (m, 10H), 7.67-7.64 (m, 3H), 7.54-7.51 (m, 1H), 7.48-7.40 (m, 6H), 7.38-7.21 (m, 7H), 7.09-7.07 (m, 1H), 7.01-6.99 (m, 2H), 6.95-6.94 (dd, 1H), 6.87-6.83 (m, 2H)

Synthesis Example 41: Synthesis of Compound 66

[0190] 5.38 g of Compound 66 (Yield 77%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-31 instead of Intermediate 4-25 was used. Compound 66 was identified using MS/FAB and ^{1}H NMR.
$C_{53}H_{34}N_2$: calc. 698.27, and found; 698.28
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.06-8.04 (dd, 1H), 7.88-7.85 (m, 1H), 7.84-7.79 (m, 4H), 7.78-7.76 (m, 1H), 7.74-7.69 (m, 4H), 7.67-7.64 (m, 2H), 7.56-7.52 (m, 1H), 7.48-7.29 (m, 13H), 7.22 (t, 1H), 7.13-7.11 (m, 1H), 7.05-7.01 (m, 2H), 6.96-6.94 (dd, 1H), 6.87-6.84 (m, 2H)

Synthesis Example 42: Synthesis of Compound 67

[0191] 5.36 g of Compound 67 (Yield 74%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-67 instead of Intermediate 4-25 was used. Compound 67 was identified using MS/FAB and ^{1}H NMR.
$C_{55}H_{36}N_2$: calc. 724.29, and found; 724.30
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.85-7.82 (m, 2H), 7.81-7.79 (m, 2H), 7.78-7.76 (m, 1H), 7.73-7.68 (m, 4H), 7.67 (d, 1H), 7.65 (d, 1H), 7.60-7.57 (m, 2H), 7.54-7.52 (dd, 1H), 7.48-7.28 (m, 13H), 7.26-7.22 (m, 2H), 7.16-7.15 (m, 1H), 7.06-7.01 (m, 2H), 6.94-6.90 (m, 2H), 6.87-6.84 (m, 2H)

Synthesis Example 43: Synthesis of Compound 68

[0192] 5.45 g of Compound 68 (Yield 78%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-68 instead of Intermediate 4-25 was used. Compound 68 was identified using MS/FAB and ^{1}H NMR.

$C_{53}H_{34}N_2$· calc. 698.27, and found; 698.28

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.86-7.83 (m, 2H), 7.82-7.79 (m, 2H), 7.78-7.76 (m, 2H), 7.74-7.68 (m, 4H), 7.66-7.63 (m, 3H), 7.57-7.54 (m, 2H), 7.53-7.50 (m, 2H), 7.47-7.43 (m, 3H), 7.41-7.28 (m, 8H), 7.18-7.16 (m, 1H), 7.06-7.05 (dd, 1H), 6.92-6.88 (m, 2H), 6.85-6.82 (m, 2H)

Synthesis Example 44: Synthesis of Compound 70

[0193] 3.56 g of Compound 70 (Yield 53%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-70 instead of Intermediate 4-25 was used. Compound 70 was identified using MS/FAB and ^{1}H NMR.

$C_{50}H_{29}N_3$: calc. 671.24, and found; 671.23

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.22-8.20 (m, 2H), 7.92-7.88 (m, 2H), 7.85-7.82 (m, 2H), 7.80 (d, 1H), 7.75-7.69 (m, 4H), 7.67-7.66 (dd, 1H), 7.65-7.63 (m, 2H), 7.61 (d, 1H), 7.55-7.53 (dd, 1H), 7.49-7.42 (m, 7H), 7.38-7.29 (m, 5H), 7.02-6.97 (m, 1H)

Synthesis Example 45: Synthesis of Compound 73

[0194] 4.97g of Compound 73 (Yield 65%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-43 instead of Intermediate 4-25 was used. Compound 73 was identified using MS/FAB and ^{1}H NMR.

$C_{58}H_{40}N_2$: calc. 764.32, and found; 764.33

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 7.85-7.82 (m, 2H), 7.81-7.79 (m, 2H), 7.78-7.75 (m, 2H), 7.73-7.68 (m, 4H), 7.67-7.64 (m, 2H), 7.56-7.52 (m, 2H), 7.47-7.44 (m, 3H), 7.39-7.27 (m, 8H), 7.13-7.09 (m, 2H), 7.02-7.00 (m, 1H), 6.96-6.92 (m, 2H), 6.86-6.84 (dd, 1H), 6.82-6.79 (m, 2H), 6.77 (d, 1H), 1.65 (s, 6H)

Synthesis Example 46: Synthesis of Compound 76

[0195] 6.06 g of Compound 76 (Yield 78%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-76 instead of Intermediate 4-25 was used. Compound 76 was identified using MS/FAB and ^{1}H NMR.

$C_{59}H_{40}N_2$: calc .776.32, and found; 776.32

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.68-8.67 (m, 1H), 8.39-8.36 (m, 1H), 7.94-7.91 (m, 1H), 7.84-7.80 (m, 3H), 7.79-7.78 (m, 1H), 7.73-7.69 (m, 4H), 7.67-7.61 (m, 4H), 7.55-7.38 (m, 10H), 7.37-7.24 (m, 7H), 7.18-7.16 (m, 1H), 7.06-7.02 (m, 2H), 6.91-6.88 (m, 2H), 6.82-6.78 (m, 2H)

Synthesis Example 47: Synthesis of Compound 77

[0196] 5.22 g of Compound 77 (Yield 72%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-77 instead of Intermediate 4-25 was used. Compound 77 was identified using MS/FAB and ^{1}H NMR.

$C_{54}H_{35}N_3$: calc. 725.28, and found; 725.27

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.55-8.52 (m, 2H), 7.85-7.82 (m, 2H), 7.81-7.77 (m, 3H), 7.74-7.70 (m, 4H), 7.67 (d, 1H), 7.65 (d, 1H), 7.58-7.52 (m, 5H), 7.47-7.44 (m, 3H), 7.39-7.28 (m, 7H), 7.11-7.09 (m, 1H), 7.02-6.99 (m, 2H), 6.93-6.89 (m, 2H), 6.81-6.77 (m, 2H)

Synthesis Example 48: Synthesis of Compound 78

[0197] 5.63 g of Compound 78 (Yield 75%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-78 instead of Intermediate 4-25 was used. Compound 78 was identified using MS/FAB and ^{1}H NMR.

$C_{57}H_{38}N_2$: calc. 750.30, and found; 750.29

^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.67-8.65 (dd, 1H), 8.40-8.37 (m, 1H), 7.95-7.92 (m, 1H), 7.85-7.82 (m, 2H), 7.81-7.79 (m, 2H), 7.78-7.76 (m, 2H), 7.74-7.68 (m, 4H), 7.66-7.63 (m, 3H), 7.58-7.52 (m, 4H), 7.48-7.44 (m, 4H), 7.37-7.25 (m,

8H), 7.16-7.15 (m, 1H), 7.02-7.00 (dd, 1H), 6.95-6.91 (m, 2H), 6.84-6.81 (m, 2H)

Synthesis Example 49: Synthesis of Compound 79

[0198] 5.00 g of Compound 79 (Yield 69%) was prepared in the same manner as in the method of preparing Compound 58 of Synthesis Example 34, except that Intermediate 4-79 instead of Intermediate 4-25 was used. Compound 79 was identified using MS/FAB and ^{1}H NMR.
$C_{54}H_{35}N_3$: calc. 725.28, and found; 725.27
^{1}H NMR ($CDCl_3$, 400Hz) δ (ppm) 8.49-8.46 (m, 1H), 7.84-7.77 (m, 7H), 7.74-7.68 (m, 5H), 7.67-7.63 (m, 3H), 7.56-7.54 (dd, 1H), 7.47-7.43 (m, 3H), 7.39-7.26 (m, 8H), 7.19-7.17 (m, 1H), 7.03-6.99 (m, 2H), 6.92-6.88 (m, 2H), 6.81-6.78 (m, 2H)

Synthesis Example 50: Synthesis of Compound 82

Synthesis of Intermediate 3-82

[0199] Intermediate 3-82 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-82 instead of Intermediate 2-1 was used.

Synthesis of Compound 82

[0200] 5.20 g of Compound 82 (Yield 67%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-82 and 4-25, instead of Intermediates 3-1 and 4-1, were used. Compound 82 was identified using MS/FAB and ^{1}H NMR.
$C_{59}H_{40}N_2$: calc. 776.32, and found; 776.33
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.68 (d, 1H), 8.41-8.38 (m, 1H), 7.94-7.91 (m, 1H), 7.86-7.80 (m, 5H), 7.74-7.64 (m, 10H), 7.56-7.53 (m, 1H), 7.49-7.44 (m, 4H), 7.38-7.24 (m, 7H), 7.13-7.09 (m, 2H), 7.04-6.99 (m, 3H), 6.96-6.92 (m, 2H), 6.87-6.83 (m, 1H), 6.81-6.78 (m, 2H)

Synthesis Example 51: Synthesis of Compound 83

[0201] 5.58 g of Compound 83 (Yield 68%) was prepared in the same manner as in the method of preparing Compound 82 of Synthesis Example 50, except that Intermediate 4-48 instead of Intermediate 4-25 was used. Compound 83 was identified using MS/FAB and ^{1}H NMR.
$C_{63}H_{42}N_2$: calc. 826.33, and found; 826.32
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.46-8.43 (m, 1H), 8.17-8.14 (m, 1H), 7.87-7.80 (m, 6H), 7.79-7.75 (m, 2H), 7.74-7.63 (m, 12H), 7.56-7.54 (dd, 1H), 7.49-7.40 (m, 6H), 7.38-7.22 (m, 7H), 7.16-7.14 (m, 1H), 7.05-7.00 (m, 2H), 6.95-6.93 (dd, 1H), 6.85-6.82 (m, 2H)

Synthesis Example 52: Synthesis of Compound 84

Synthesis of Intermediate 3-84

[0202] Intermediate 3-84 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-84 instead of Intermediate 2-1 was used.

Synthesis of Compound 84

[0203] 4.90 g of Compound 84 (Yield 62%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-84 and 4-30, instead of Intermediates 3-1 and 4-1, were used. Compound 84 was identified using MS/FAB and ^{1}H NMR.
$C_{60}H_4{=}N_2$: calc. 790.33, and found; 790.32
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.69 (d, 1H), 8.40-8.37(m, 1H), 8.11-8.08 (m, 1H), 7.93-7.90 (m, 2H), 7.89-7.85 (m, 2H), 7.84-7.79 (m, 4H), 7.72-7.68 (m, 2H), 7.62-7.59 (m, 2H), 7.53-7.51 (m, 1H), 7.49-7.42 (m, 6H), 7.40-7.25 (m, 8H), 7.13-7.11 (m, 1H), 7.07-7.05 (dd, 1H), 6.98-6.96 (dd, 1H), 6.90-6.86 (m, 2H), 6.82 (d, 1H), 1.62 (s, 6H)

Synthesis Example 53 : Synthesis of Compound 85

[0204] 4.90 g of Compound 84 (Yield 64%) was prepared in the same manner as in the method of preparing Compound

84 of Synthesis Example 52, except that Intermediate 4-27 instead of Intermediate 4-30 was used. Compound 85 was identified using MS/FAB and ^{1}H NMR.
$C_{57}H_{39}N_3$: calc. 765.31, and found; 765.30
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.68 (d, 1H), 8.38-8.35 (m, 1H), 7.95-7.92 (m, 2H), 7.90-7.88 (m, 1H), 7.85-7.80 (m, 4H), 7.73-7.69 (m, 2H), 7.63-7.60 (m, 2H), 7.55-7.52 (m, 2H), 7.48-7.44 (m, 2H), 7.41-7.33 (m, 5H), 7.32-7.25 (m, 4H), 7.12-7.09 (m, 1H), 7.03-6.96 (m, 2H), 6.93-6.90 (dd, 1H), 6.88-6.84 (m, 2H), 6.80 (d, 1H), 1.61 (s, 6H)

Synthesis Example 54: Synthesis of Compound 86

[0205] 4.73 g of Compound 86 (Yield 64%) was prepared in the same manner as in the method of preparing Compound 84 of Synthesis Example 52, except that Intermediate 4-31 instead of Intermediate 4-30 was used. Compound 86 was identified using MS/FAB and ^{1}H NMR.
$C_{56}H_{38}N_2$: calc. 738.30, and found; 738.31
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.11-8.08 (m, 1H), 7.94-7.91 (m, 1H), 7.89-7.85 (m, 2H), 7.84-7.79 (m, 4H), 7.72-7.69 (m, 2H), 7.62-7.59 (m, 2H), 7.54-7.51 (m, 1H), 7.49-7.41 (m, 5H), 7.40-7.34 (m, 5H), 7.32-7.25 (m, 3H), 7.14-7.12 (m, 1H), 7.01-6.99 (m, 1H), 6.93-6.87 (m, 3H), 6.85 (d, 1H), 1.62 (s, 6H)

Synthesis Example 55: Synthesis of Compound 89

Synthesis of Intermediate 3-89

[0206] Intermediate 3-89 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-89 instead of Intermediate 2-1 was used.

Synthesis of Compound 89

[0207] 5.61 g of Compound 89 (Yield 75%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-89 and 4-4, instead of Intermediates 3-1 and 4-1, were used. Compound 89 was identified using MS/FAB and ^{1}H NMR.
$C_{57}H_{36}N_2$: calc. 748.29, and found; 748.30
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.25-8.19 (m, 2H), 8.11-8.09 (m, 1H), 8.06-8.04 (m, 1H), 7.85-7.82 (m, 2H), 7.81-7.80 (m, 1H), 7.78-7.77 (m, 1H), 7.72-7.68 (m, 3H), 7.64-7.61 (m, 2H), 7.59-7.55 (m, 3H), 7.54-7.48 (m, 3H), 7.47-7.43 (m, 3H), 7.42-7.35 (m, 5H), 7.34-7.27 (m, 3H), 7.12-7.08 (m, 2H), 7.02-6.98 (m, 2H), 6.90-6.86 (m, 2H)

Synthesis Example 56: Synthesis of Compound 92

Synthesis of Intermediate 3-92

[0208] Intermediate 3-92 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-92 instead of Intermediate 2-1 was used.

Synthesis of Compound 92

[0209] 4.61 g of Compound 92 (Yield 61%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-92 and 4-27, instead of Intermediates 3-1 and 4-1, were used. Compound 92 was identified using MS/FAB and ^{1}H NMR.
$C_{54}H_{33}N_3S$: calc. 755.24, and found; 755.25
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.70-8.68 (m, 1H), 8.40-8.37 (m, 2H), 8.01 (d, 1H), 7.93-7.91 (m, 1H), 7.88-7.86 (dd, 1H), 7.85-7.82 (m, 2H), 7.81-7.80 (m, 1H), 7.78-7.76 (m, 1H), 7.73-7.69 (m, 2H), 7.64 (d, 1H) 7.63-7.62 (m, 1H), 7.54-7.51 (m, 1H), 7.48-7.43 (m, 2H), 7.40-7.35 (m, 4H), 7.34-7.24 (m, 5H), 7.13-7.12 (m, 1H), 7.06-7.03 (dd, 1H), 6.98-6.93 (m, 1H), 6.92-6.84 (m, 4H)

Synthesis Example 57: Synthesis of Compound 95

Synthesis of Intermediate 3-95

[0210] Intermediate 3-95 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-95 instead of Intermediate 2-1 was used.

Synthesis of Compound 95

[0211] 4.74 g of Compound 95 (Yield 62%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-95 and 4-30, instead of Intermediates 3-1 and 4-1, were used. Compound 95 was identified using MS/FAB and ^{1}H NMR.
$C_{57}H_{36}N_2O$: calc. 764.28, and found; 764.29
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.69-8.67 (m, 1H), 8.41-8.38 (m, 1H), 8.11-8.08 (m, 1H), 8.04-8.02 (m, 1H), 7.99-7.96 (m, 2H), 7.94-7.91 (m, 1H), 7.87-7.85 (m, 1H), 7.84-7.77 (m, 6H), 7.73-7.69 (m, 2H), 7.55-7.52 (m, 1H), 7.48-7.41 (m, 5H), 7.40-7.36 (m, 3H), 7.35-7.25 (m, 5H), 7.10-7.04 (m, 2H), 6.98-6.96 (m, 1H), 6.94-6.92 (dd, 1H), 6.88-6.85 (m, 2H)

Synthesis Example 58: Synthesis of Compound 97

Synthesis of Intermediate 3-97

[0212] Intermediate 3-97 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-97 instead of Intermediate 2-1 was used.

Synthesis of Compound 97

[0213] 4.96 g of Compound 97 (Yield 61%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-97 and 4-4, instead of Intermediates 3-1 and 4-1, were used. Compound 97 was identified using MS/FAB and ^{1}H NMR.
$C_{61}H_{39}N_3$: calc. 813.31, and found; 813.32
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.31-8.29 (m, 1H), 7.84-7.82 (m, 2H), 7.81-7.80 (m, 1H), 7.77-7.74 (m, 1H), 7.72-7.68 (m, 4H), 7.64-7.61 (m, 2H), 7.60-7.58 (m, 1H), 7.54-7.47 (m, 7H), 7.46-7.41 (m, 4H), 7.40-7.36 (m, 5H), 7.34-7.28 (m, 4H), 7.24-7.22 (m, 1H), 7.16-7.14 (m, 1H), 7.06-7.00 (m, 2H), 6.86-6.80 (m, 3H)

Synthesis Example 59: Synthesis of Compound 98

[0214] 4.65 g of Compound 98 (Yield 59%) was prepared in the same manner as in the method of preparing Compound 97 of Synthesis Example 58, except that Intermediate 4-31 instead of Intermediate 4-4 was used. Compound 98 was identified using MS/FAB and ^{1}H NMR.
$C_{59}H_{37}N_3$: calc. 787.30, and found; 787.29
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.32-8.30 (m, 1H), 8.10-8.07 (dd, 1H), 7.87-7.85 (m, 1H), 7.84-7.80 (m, 2H), 7.80-7.79 (m, 1H), 7.76-7.74 (m, 1H), 7.73-7.68 (m, 4H), 7.61-7.58 (m, 1H), 7.55-7.51 (m, 1H), 7.50-7.43 (m, 7H), 7.42-7.35 (m, 6H), 7.34-7.23 (m, 6H), 7.13-7.10 (m, 1H), 7.04-7.03 (dd, 1H), 6.94-6.91 (dd, 1H), 6.87-6.84 (m, 2H)

Synthesis Example 60: Synthesis of Compound 99

Synthesis of Intermediate 3-99

[0215] Intermediate 3-99 was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 2-98 instead of Intermediate 2-1 was used.

Synthesis of Compound 99

[0216] 5.03 g of Compound 99 (Yield 55%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-99 and 4-30, instead of Intermediates 3-1 and 4-1, were used. Compound 99 was identified using MS/FAB and ^{1}H NMR.
$C_{69}H_{45}N_3$: calc. 915.36, and found; 915.35
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.90 (s, 1H), 8.60-8.58 (dd, 1H), 8.45 (s, 1H), 8.13-8.11 (dd, 1H), 7.94-7.91 (m, 1H), 7.87-7.75 (m, 5H), 7.72-7.58 (m, 7H), 7.54-7.21 (m, 23), 7.16-7.14 (m, 1H), 7.03-6.99 (m, 2H), 6.89-6.84 (m, 2H)

Synthesis Example 61: Synthesis of Compound 103

[0217] 5.71 g of Compound 103 (Yield 76%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate B37 instead of Intermediate 4-1 was used. Compound 103 was identified using MS/FAB and ^{1}H NMR.

$C_{56}H_{37}N_3$: calc. 751.30, found 751.28
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.76 (s, 2H), 8.57 (d, 2H), 8.12 (d, 1H), 7.93 (d, 2H), 7.87-7.65 (m, 9H), 7.61-7.27 (m, 15H), 7.14-7.10 (m, 2H), 7.08-7.05 (m, 3H), 6.98 (d, 1H)

Synthesis Example 62: Synthesis of Compound 104

[0218] 5.23 g of Compound 104 (Yield 72%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediate B38 instead of Intermediate 4-1 was used. Compound 104 was identified using MS/FAB and ^{1}H NMR.
$C_{53}H_{34}N_4$: calc. 726.28, found 726.27
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.78 (s, 2H), 8.56 (d, 2H), 7.94 (d, 2H), 7.84-7.80 (m, 3H), 7.72-7.60 (m, 7H), 7.54-7.25 (m, 11H), 7.12-7.08 (m, 5H), 7.02-6.99 (m, 2H)

Synthesis Example 63: Synthesis of Compound 107

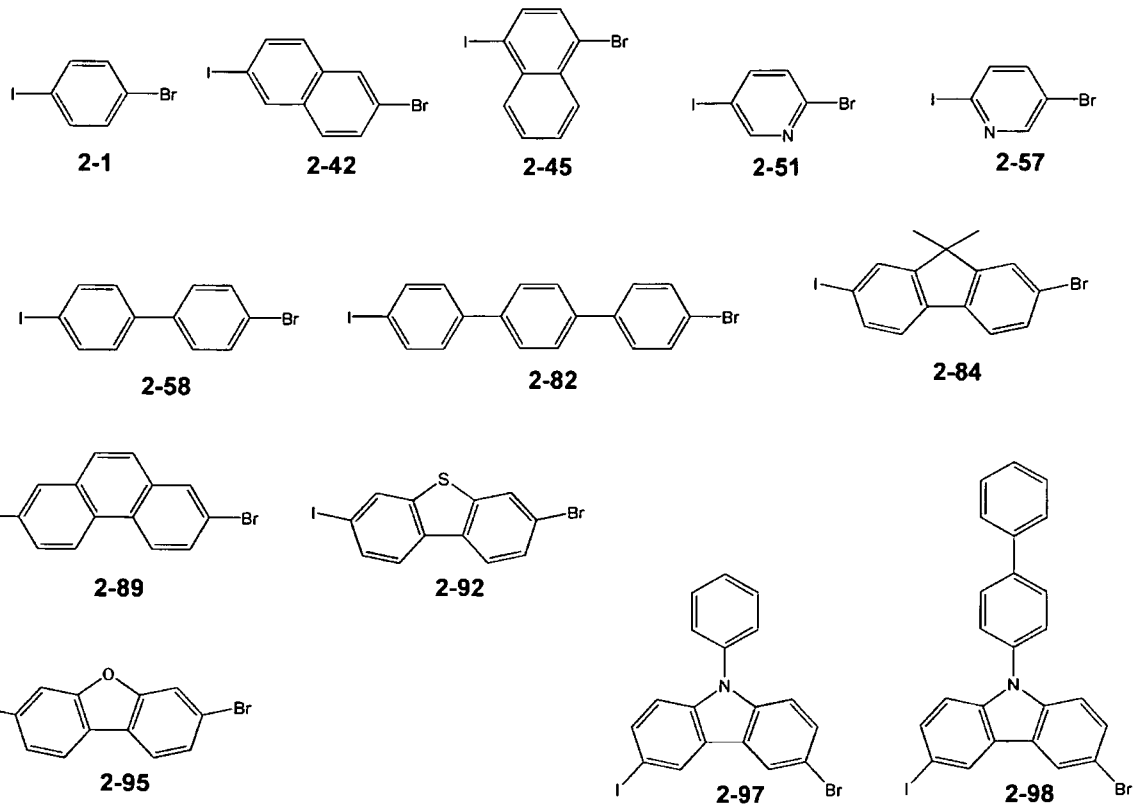
[0219] 6.21 g of Compound 107 (yield 75%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-58 and B37, instead of Intermediates 3-1 and 4-1, were used. Compound 107 was identified using MS/FAB and ^{1}H NMR.

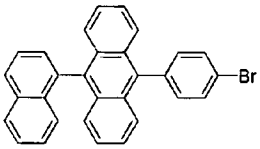
Synthesis of Compound 107

[0220] 6.21g of Compound 107 (Yield 75%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that Intermediates 3-107 and B37, instead of Intermediates 3-1 and 4-1, were used. Compound 107 was identified using MS/FAB and ^{1}H NMR.
$C_{62}H_{41}N_3$: calc. 827.33, found 827.31
^{1}H NMR ($CDCl_3$, 400MHz) δ (ppm) 8.76 (s, 2H), 8.57 (d, 2H), 8.11 (d, 1H), 7.94 (d, 2H), 7.87-7.62 (m, 13H), 7.54-7.26 (m, 15H), 7.12-7.10 (m, 3H), 7.06-7.03 (m, 1H), 6.96-6.92 (m, 2H)

Intermediates

[0221]





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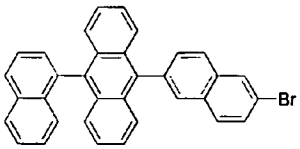

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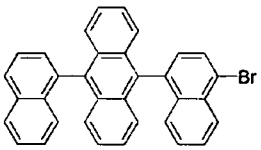

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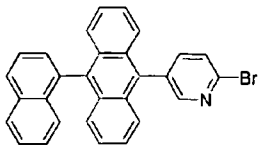

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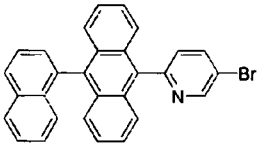

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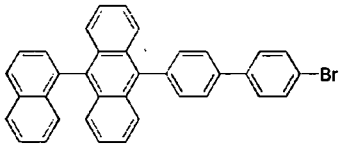

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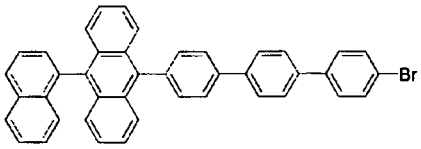

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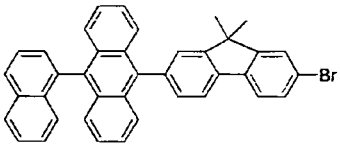

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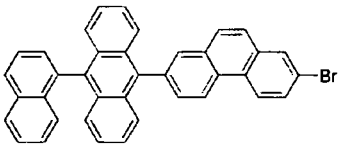

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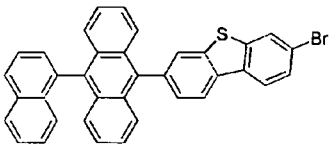

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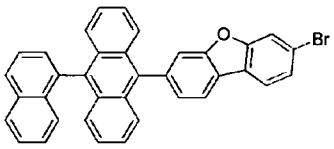

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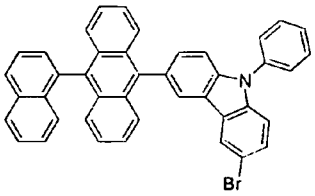

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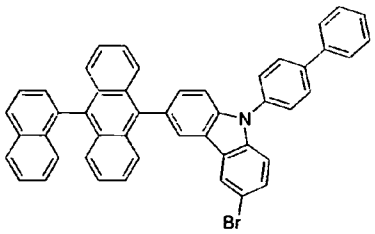

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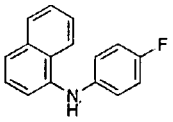

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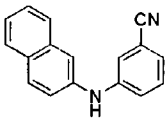

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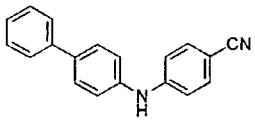

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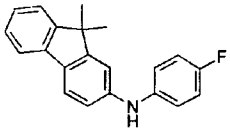

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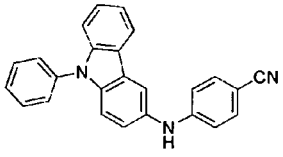

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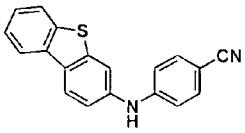

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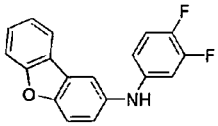

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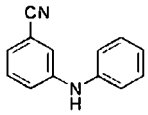

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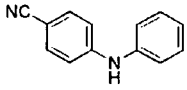

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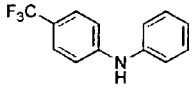

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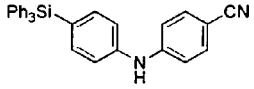

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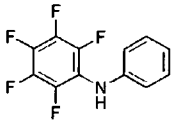

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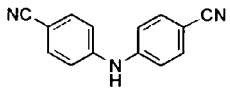

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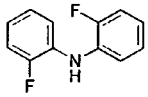

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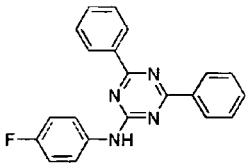

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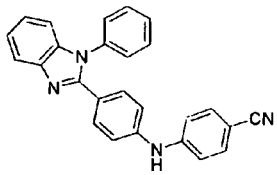

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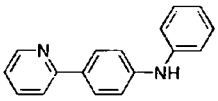

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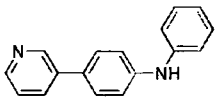

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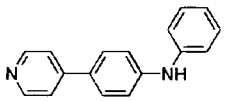

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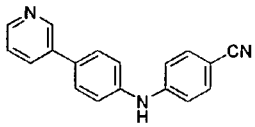

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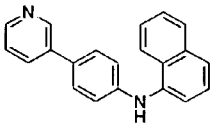

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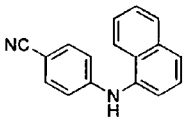

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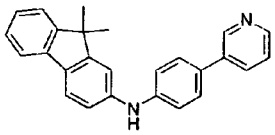

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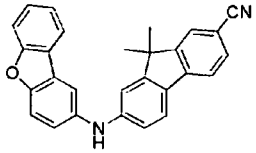

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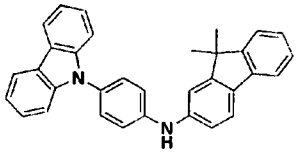

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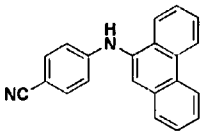

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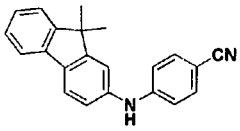

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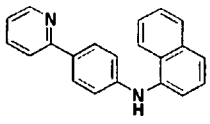

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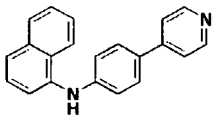

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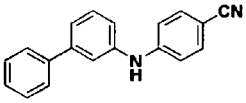

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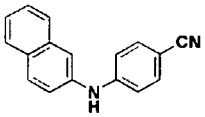

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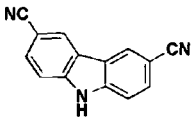

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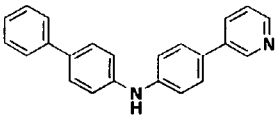

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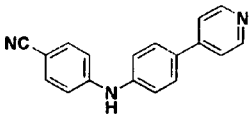

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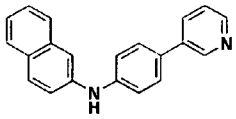

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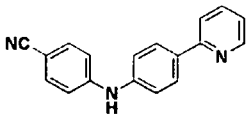

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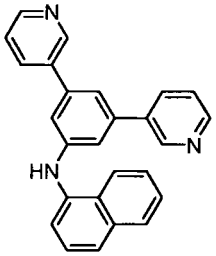

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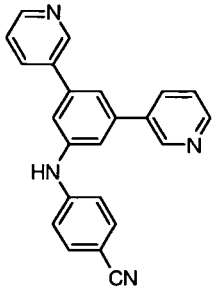


B38

Example 1

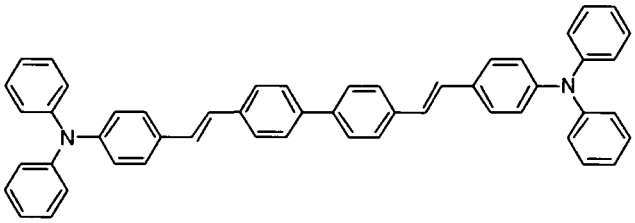
[0222] To manufacture an anode, a coming 15 Ω/cm^2 (1200 Å) ITO glass substrate was cut to a size of 50 mm x 50 mm x 0.7 mm and then sonicated in isopropyl alcohol and pure water each for five minutes, and then cleaned by irradiation of ultraviolet rays for 30 minutes and exposure to ozone. The resulting glass substrate was loaded into a vacuum deposition diode.

[0223] 2-TNATA was deposited on the ITO glass substrate to form an HIL having a thickness of 600Å on the anode, and then 4,4'-bis[N-(1-naphthyl)-N-phenylamino]biphenyl (NPS) was deposited on the HIL to form a HTL having a thickness of 300Å.

[0224] Subsequently, 9,10-di-naphthalene-2-yl-anthracene (ADN) and 4,4'-bis[2-(4-(N,N-diphenylamino)phenyl)vinyl]biphenyl (DPAVBi) were co-deposited on the HTL in a weight ratio of 98:2 to form an EML having a thickness of about 300Å.

[0225] Then, Compound 1 was deposited on the EML to form an ETL having a thickness of about 300Å, and then LiF was deposited on the ETL to form an EIL having a thickness of about 10Å. Then, Al was deposited on the EIL to form a second electrode (cathode) having a thickness of about 3000Å, thereby completing the manufacture of an organic light-emitting diode.

<DPAVBi>



Example 2

[0226] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 4, instead of Compound 1, was used to form the ETL.

Example 1

[0227] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 14, instead of Compound 1, was used to form the ETL.

Example 4

[0228] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 23, instead of Compound 1, was used to form the ETL.

Example 5

[0229] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 25, instead of Compound 1, was used to form the ETL.

Example 6

[0230] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 27, instead of Compound 1, was used to form the ETL.

Example 7

[0231] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 31, instead of Compound 1, was used to form the ETL.

Example 8

[0232] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 32, instead of Compound 1, was used to form the ETL.

Example 9

[0233] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 42, instead of Compound 1, was used to form the ETL.

Example 10

[0234] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 48, instead of Compound 1, was used to form the ETL.

Example 11

[0235] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 58, instead of Compound 1, was used to form the ETL.

Example 12

[0236] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 60, instead of Compound 1, was used to form the ETL.

Example 13

[0237] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 62, instead of Compound 1, was used to form the ETL.

Example 14

[0238] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 66, instead of Compound 1, was used to form the ETL.

Example 15

[0239] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 70, instead of Compound 1, was used to form the ETL.

Example 16

[0240] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 77, instead of Compound 1, was used to form the ETL.

Example 17

[0241] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 82, instead of Compound 1, was used to form the EML.

Example 18

[0242] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 86, instead of Compound 1, was used to form the ETL.

Example 19

[0243] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound

97, instead of Compound 1, was used to form the EML.

Example 20

[0244] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 25, instead of DPAVBi, was used to form the EML, and Alq_3, instead of Compound 1, was used to form the ETL.

Example 21

[0245] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound 86, instead of DPAVBi, was used to form the EML, and Alq_3, instead of Compound 1, was used to form the ETL.

Example 22

[0246] An organic light-emitting device was manufactured in the same manner as in Example 1, except that Compound 103, instead of Compound 1, was used to form the ETL.

Example 23

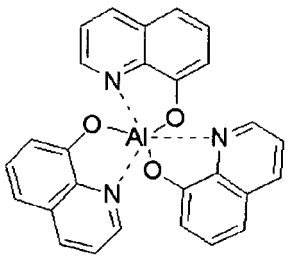
[0247] An organic light-emitting device was manufactured in the same manner as in Example 1, except that Compound 104, instead of Compound 1, was used to form the ETL.

Example 24

[0248] An organic light-emitting device was manufactured in the same manner as in Example 1, except that Compound 107, instead of Compound 1, was used to form the ETL.

Comparative Example 1

[0249] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Alq_3, instead of Compound 1, was used to form the ETL.

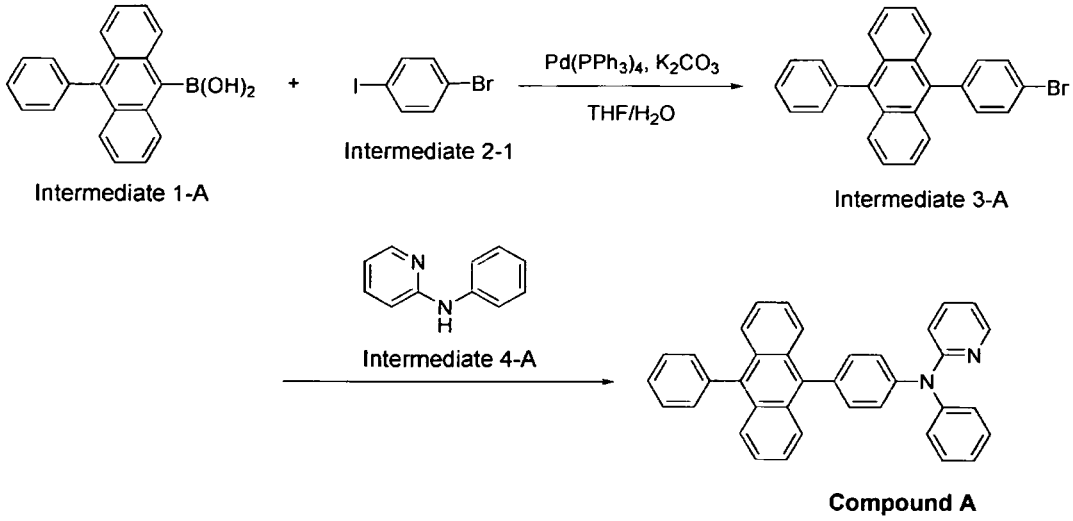


Alq_3

Comparative Example 2

[0250] Compound A was synthesized according to Reaction Scheme A below:

Reaction Scheme A



Synthesis of Intermediate 3-A

[0251] Intermediate 3-A was prepared in the same manner as in the method of preparing Intermediate 3-1 of Synthesis Example 1, except that Intermediate 1-A, instead of Intermediate 1, was used.

Synthesis of Compound A

[0252] 3.23 g of Compound A (Yield 63%) was prepared in the same manner as in the method of preparing Compound 1 of Synthesis Example 1, except that 4.24 g (10 mmol) of Intermediate 3-A and 2.23 g (12.0 mmol) of Intermediate 4-A, instead of Intermediates 3-1 and 4-1, were used.

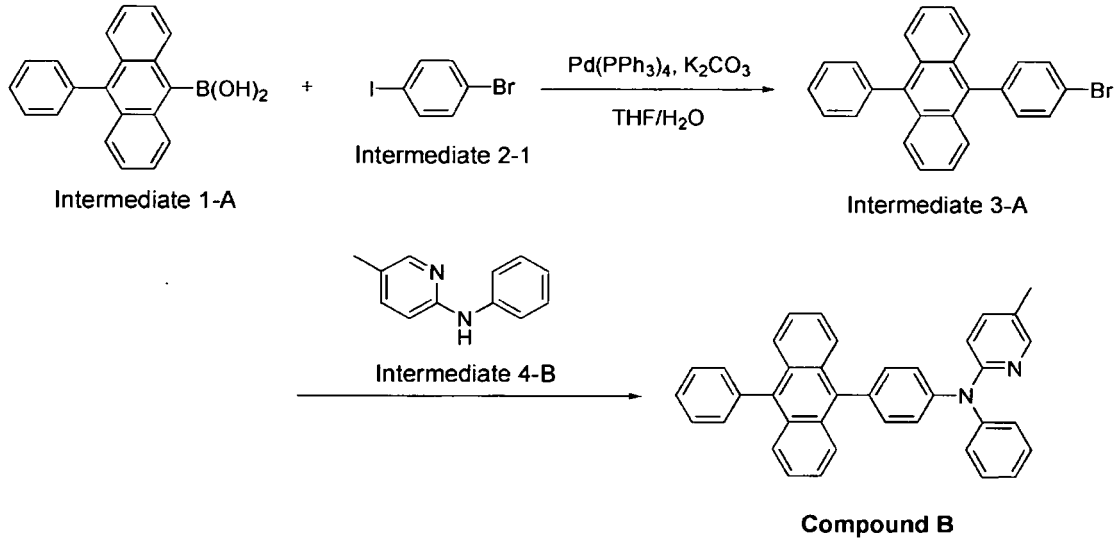
Manufacture of organic light-emitting diode

[0253] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound A, instead of Compound 1, was used to form the ETL.

Comparative Example 3

[0254] Compound B was synthesized according to Reaction Scheme B below:

Reaction Scheme B



Synthesis of Compound B

[0255] Intermediate B was prepared in the same manner as in the method of preparing Compound A of Comparative Example 2, except that Intermediate 4-B, instead of Intermediate 4-A, was used.

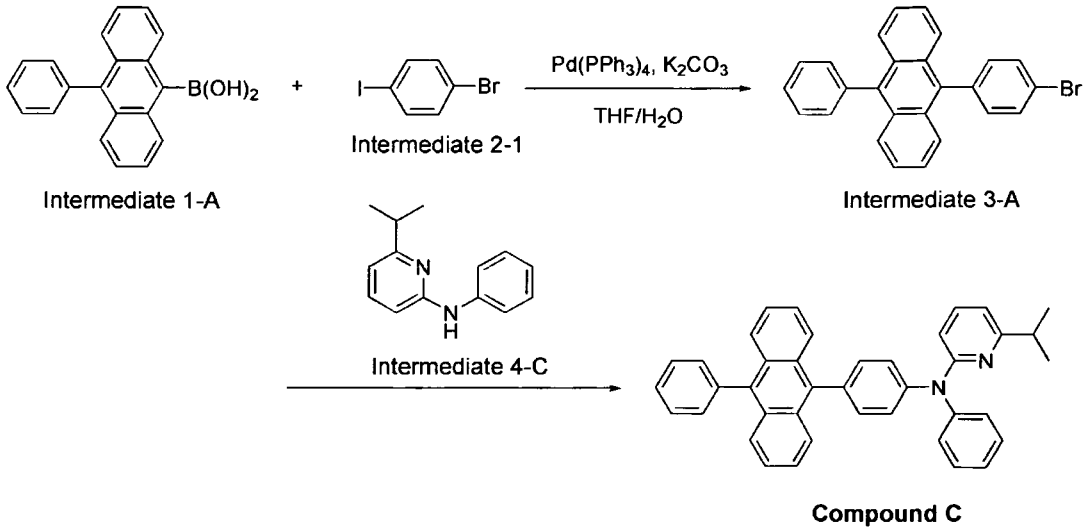
Manufacture of organic light-emitting diode

[0256] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound B, instead of Compound 1, was used to form the ETL.

Comparative Example 4

[0257] Compound C was synthesized according to Reaction Scheme C below:

Reaction Scheme C



Synthesis of Compound C

[0258] Compound C was prepared in the same manner as in the method of preparing Compound A of Comparative Example 2, except that Intermediate 4-C, instead of Intermediate 4-A, was used.

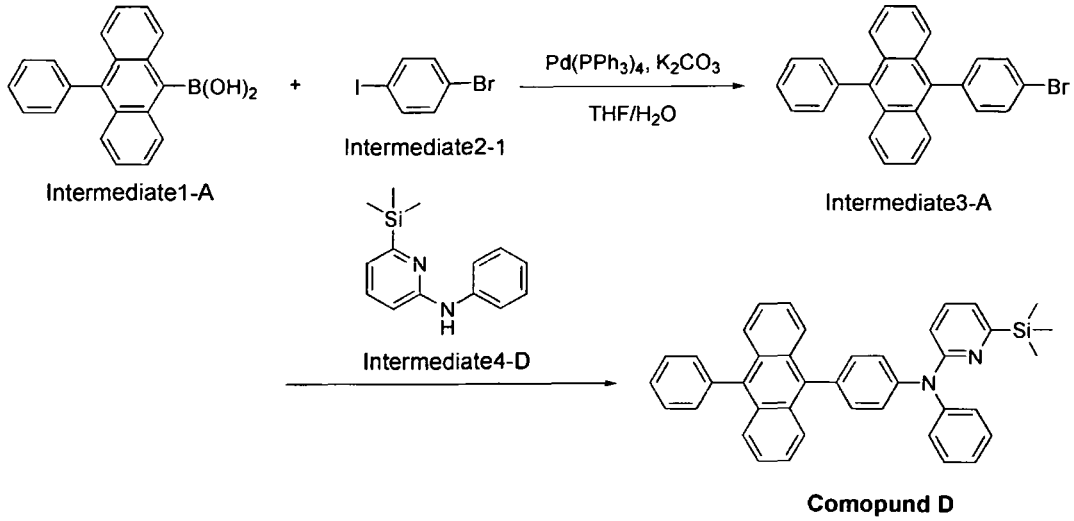
Manufacture of organic light-emitting diode

[0259] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound C, instead of Compound 1, was used to form the ETL.

Comparative Example 5

[0260] Compound D was synthesized according to Reaction Scheme D below:

Reaction Scheme D



Synthesis of Compound D

[0261] Compound D was prepared in the same manner as in the method of preparing Compound A of Comparative Example 2, except that Intermediate 4-D, instead of Intermediate 4-A, was used.

Manufacture of organic light-emitting diode

[0262] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that Compound D, instead of Compound 1, was used to form the ETL.

Evaluation Example 1

[0263] Driving voltages, luminances, emitting-light colours, efficiencies (at current density of 50mA/cm^2, and half-life spans (at 100mA/cm^2) of the organic light-emitting diodes of Examples 1 to 24 and Comparative Examples 1 to 5 were measured using a PR650 (Spectroscan) Source Measurement Unit (available from Photo Research, Inc.). The results are shown in Table 1 below.

Table 1

	EML host	EML dopant	ETL	Driving voltage (V)	Luminance (cd/m^2)	Efficiency (cd/A)	Light-emitting colour	Half-life span (hr)
Example 1	ADN	DPVABi	Compound 1	5.36	3,335	6.67	blue	492
Example 2	ADN	DPVABi	Compound 4	5.32	3,390	6.78	blue	596
Example 3	ADN	DPVABi	Compound 14	5.29	3,380	6.76	blue	542

(continued)

	EML host	EML dopant	ETL	Driving voltage (V)	Luminance (cd/m^2)	Efficiency (cd/A)	Light-emitting colour	Half-life span (hr)
Example 4	ADN	DPVABi	Compound 23	5.24	3,470	6.94	blue	589
Example 5	ADN	DPVABi	Compound 25	5.26	3,465	6.93	blue	536
Example 6	ADN	DPVABi	Compound 27	5.21	3,505	7.01	blue	564
Example 7	ADN	DPVABi	Compound 31	5.26	3,280	6.56	blue	632
Example 8	ADN	DPVABi	Compound 32	5.30	3,440	6.88	blue	559
Example 9	ADN	DPVABi	Compound 42	5.23	3,565	7.13	blue	658
Example 10	ADN	DPVABi	Compound 48	5.31	3,435	6.87	blue	524
Example 11	ADN	DPVABi	Compound 58	5.26	3,580	7.16	blue	532
Example 12	ADN	DPVABi	Compound 60	5.24	3,515	7.03	blue	582
Example 13	ADN	DPVABi	Compound 62	5.31	3,605	7.21	blue	529
Example 14	ADN	DPVABi	Compound 66	5.27	3,345	6.69	blue	608
Example 15	ADN	DPVABi	Compound 70	5.16	3,015	6.03	blue	583
Example 16	ADN	DPVABi	Compound 77	5.21	3,560	7.12	blue	637
Example 17	ADN	DPVABi	Compound 82	5.32	3,460	6.92	blue	543
Example 18	ADN	DPVABi	Compound 86	5.43	3,445	6.89	blue	469
Example 19	ADN	DPVABi	Compound 97	5.62	3,070	6.14	blue	486
Example 20	ADN	Compound 25	Alq_3	6.86	2,190	4.38	blue	216
Example 21	ADN	Compound 86	Alq_3	6.97	2,260	4.52	blue	238
Example 22	ADN	DPVABi	Compound 103	5.35	3,640	7.82	blue	659
Example 23	ADN	DPVABi	Compound 104	5.12	3,445	6.89	blue	618
Example 24	ADN	DPVABi	Compound 107	5.63	3,590	7.18	blue	689

(continued)

	EML host	EML dopant	ETL	Driving voltage (V)	Luminance (cd/m^2)	Efficiency (cd/A)	Light-emitting colour	Half-life span (hr)
Comparative Example 1	ADN	DPVABi	Alq_3	7.35	2,065	4.13	blue	145
Comparative Example 2	ADN	DPVABi	Compound A	5.71	2,865	5.73	blue	311
Comparative Example 3	ADN	DPVABi	Compound B	5.75	2,730	5.46	blue	320
Comparative Example 4	ADN	DPVABi	Compound C	5.73	2,845	5.69	blue	297
Comparative Example 5	ADN	DPVABi	Compound D	5.73	2,830	5.66	blue	213

[0264] Referring to Table 1, the organic light-emitting devices of Examples 1 to 19 and 22 to 24 are found to have lower driving voltages, higher luminances, higher efficiencies, and better lifetime characteristics as compared to the organic light-emitting diodes of Comparative Examples 1 to 5. The organic light-emitting devices of Examples 20 and 21 are found to have lower driving voltages and better lifetime characteristics as compared to the organic light-emitting diode of Comparative Example 1.

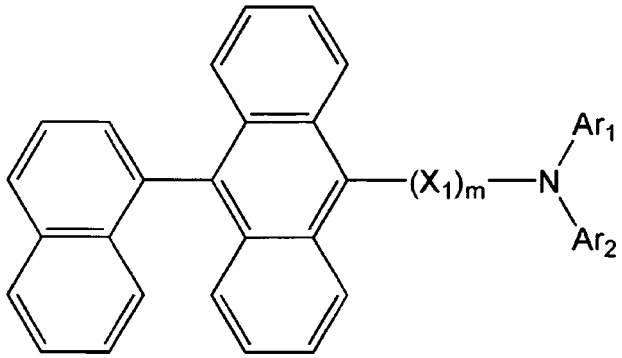
[0265] As described above, an organic light-emitting diode including any of the amine-based compounds according to the invention in its first aspect may have a low driving voltage, a high luminance, a high efficiency, and a long lifetime.

[0266] While the present invention has been described in connection with certain exemplary embodiments thereof, it is to be understood that the invention is not limited to the disclosed embodiments but, on the contrary, is intended to cover various modifications and equivalent arrangements included within the scope of the appended claims.

Claims

1. An amine-based compound represented by Formula 1 below:

Formula 1



wherein, in Formula 1, Ar_1 and Ar_2 are each independently a substituted or unsubstituted C_6-C_{60} aryl group or a substituted or unsubstituted C_2-C_{60} heteroaryl group; and wherein Ar_1 and Ar_2 are optionally linked by a single bond;

X_1 is a substituted or unsubstituted C_6-C_{60} arylene group or a substituted or unsubstituted C_2-C_{60} heteroarylene group;

m is an integer from 1 to 5; and

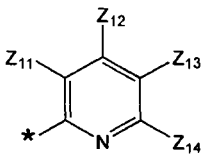
at least one substituent of each of the substituted C_6-C_{60} aryl group, the substituted C_2-C_{60} heteroaryl group, the substituted C_6-C_{60}arylene group, and the substituted C_2-C_{60} heteroarylene group is one of a deuterium

atom; -F; -Cl; -Br; -I; -CN; a hydroxyl group; $-NO_2$; an amino group; an amidino group; hydrazine; hydrazone; a carboxyl group or a salt thereof; a sulfonic acid group or a salt thereof; a phosphoric acid or a salt thereof; a tri(C_6-C_{60}aryl)silyl group; a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group; a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group and a C_2-C_{60} alkynyl group that is substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof; a C_3-C_{60} cycloalkyl group, a C_3-C_{60} cycloalkenyl group, a C_6-C_{60} aryl group, a C_2-C_{60} heteroaryl group, a C_6-C_{60} aralkyl group, a C_6-C_{60} aryloxy group, a C_6-C_{60} arylthiol group; a C_3-C_{60} cycloalkyl group, a C_3-C_{60} cycloalkenyl group, a C_6-C_{60} aryl group, a C_2-C_{60} heteroaryl group, a C_6-C_{60} aralkyl group, a C_6-C_{60} aryloxy group, and a C_6-C_{60} arylthiol group that is substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one fluorine (F), a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group and a C_2-C_{60} heteroaryl group,

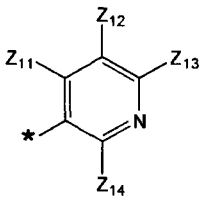
wherein at least one of Ar_1 and Ar_2 is a C_6-C_{60} aryl group substituted with at least one electron withdrawing group selected from the group consisting of -F; -CN; $-NO_2$; a C_1-C_{60} alkyl group substituted with at least one -F; a C_2-C_{60} heteroaryl group; and a C_2-C_{60} heteroaryl group substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, - CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one -F, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group, and a C_2-C_{60} heteroaryl group.

2. The amine-based compound of claim 1, wherein the at least one electron withdrawing group is selected from the group consisting of: -F; -CN; $-NO_2$; a C_1-C_{20} alkyl group substituted with at least one -F; a C_2-C_{20} heteroaryl group including a ring-containing a N atom; and a C_2-C_{20} heteroaryl group that includes a ring-containing a N atom and is substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{20}alkyl group, a C_1-C_{20} alkyl group substituted with at least one - F, a C_1-C_{20} alkoxy group, a C_6-C_{20} aryl group, and a C_2-C_{20} heteroaryl group.

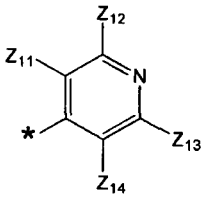
3. The amine-based compound of claim 1 or claim 2, wherein the at least one electron withdrawing group is selected from the group consisting of -F; -CN; $-CH_2F$; - CHF_2; $-CF_3$; and groups represented by Formulae 2(1) to 2(14) below:



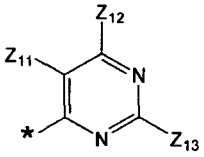
Formula 2(1)



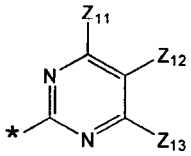
Formula 2(2)



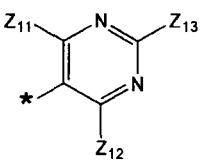
Formula 2(3)



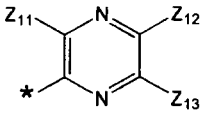
Formula 2(4)



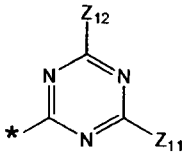
Formula 2(5)



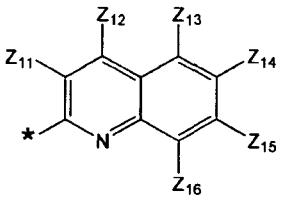
Formula 2(6)



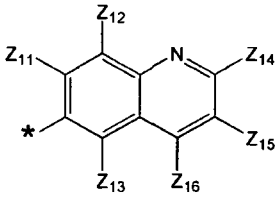
Formula 2(7)



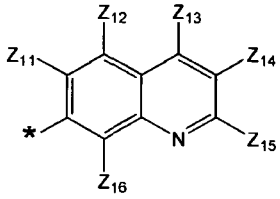
Formula 2(8)



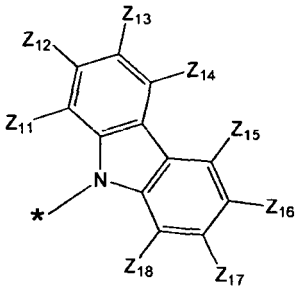
Formula 2(9)



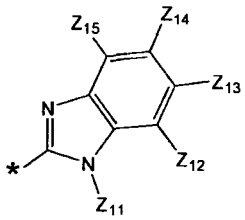
Formula 2(10)



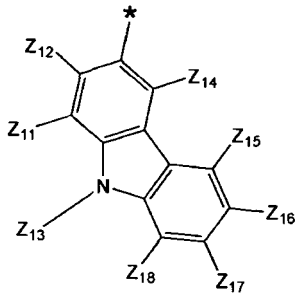
Formula 2(11)



Formula 2(12)



Formula 2(13)



Formula 2(14)

wherein, in Formulae 2(1) to 2(14), Z_{11} to Z_{18} are each independently a hydrogen atom, a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, or a carbazolyl group.

4. The amine-based compound of any preceding claim, wherein Ar_1 and Ar_2 are each independently a C_6-C_{60} aryl group selected from a substituted or unsubstituted phenyl group, a substituted or unsubstituted naphthyl group, a substituted or unsubstituted fluorenyl group, a substituted or unsubstituted phenanthrenyl group, a substituted or unsubstituted anthryl group, a substituted or unsubstituted triphenylenyl group, a substituted or unsubstituted pyrenyl group, a substituted or unsubstituted chrysenyl group, a substituted or unsubstituted pyridinyl group, a substituted or unsubstituted pyrazinyl group, a substituted or unsubstituted pyrimidinyl group, a substituted or unsubstituted quinolyl group, a substituted or unsubstituted carbazolyl group, a substituted or unsubstituted triazinyl group, a substituted or unsubstituted dibenzothiophenyl group, a substituted or unsubstituted dibenzofuranyl group, or a substituted or unsubstituted phenanthrolinyl group.

5. The amine-based compound of any preceding claim, wherein Ar_1 and Ar_2 are linked by a single bond.

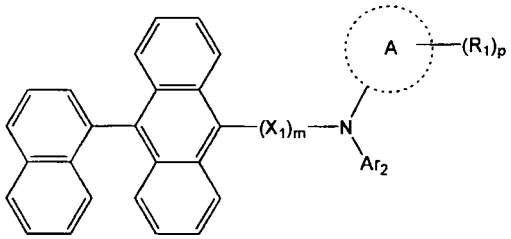
6. The amine-based compound of any preceding claim, wherein the at least one of Ar_1 and Ar_2 is a C_6-C_{60} aryl group substituted with at least two electron withdrawing groups.

7. The amine-based compound of any preceding claim, wherein the at least one of Ar_1 and Ar_2 is a phenyl group, a biphenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, or a fluorenyl group that is substituted with at least two electron withdrawing groups; and
the electron withdrawing groups are each independently selected from the group consisting of a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; and a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a phthalazinyl group, a benzoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN; a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group.

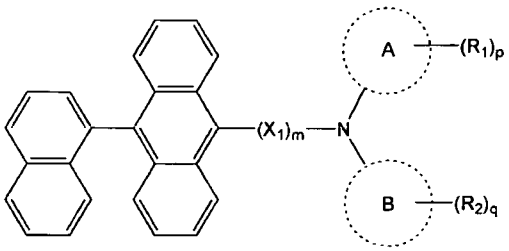
8. The amine-based compound of claim 1, wherein the amine-based compound is represented by Formula 1(1) or

1(2) below:

Formula 1(1)



Formula 1(2)



wherein, in Formula 1(1), Ar_2 is a substituted or unsubstituted C_6-C_{20} aryl group or a substituted or unsubstituted C_2-C_{20} heteroaryl group; and
in Formulae 1(1) and 1(2), the ring groups A and B are each independently a substituted C_6-C_{20} aryl group;
R_1 and R_2 are each independently an electron withdrawing group selected from the group consisting of -F; -CN; $-NO_2$; a C_1-C_{60} alkyl group substituted with at least one -F; a C_2-C_{60} heteroaryl group; and a C_2-C_{60} heteroaryl group substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one -F, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group, and a C_2-C_{60}heteroaryl group; and
p and q are each independently an integer from 1 to 9.

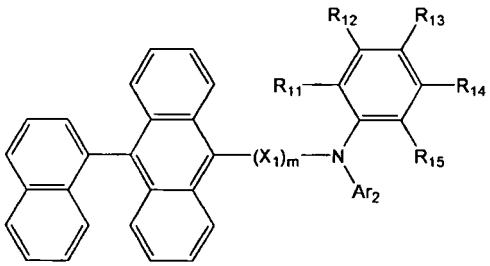
9. The amine-based compound of claim 8, wherein the amine-based compound is represented by Formula 1(1) in which at least one of R_1 is -CN; or is represented by Formula 1(2) in which at least one of R_1 and one of R_2 is -CN.

10. The amine-based compound of claim 7, wherein the amine-based compound is represented by Formula 1(1), wherein the ring group A is a substituted phenyl group, a substituted biphenyl group, a substituted naphthyl group, a substituted anthryl group, a substituted phenanthrenyl group, a substituted pyrenyl group, or a substituted fluorenyl group;
R_1 is at least one electron withdrawing group selected from the group consisting of a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; and a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a phthalazinyl group, a benzoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN; a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group; and p is 2, 3, or 4; or., wherein the amine-based compound is represented by Formula 1(2), wherein the ring group A and the ring group B are each independently a substituted phenyl group, a substituted biphenyl group, a substituted naphthyl group, a substituted anthryl group, a substituted phenanthrenyl group, a substituted pyrenyl group, or a substituted fluorenyl group;
R_1 and R_2 are each independently at least one electron withdrawing group selected from the group consisting of a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; and a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a phthalazinyl group, a ben-

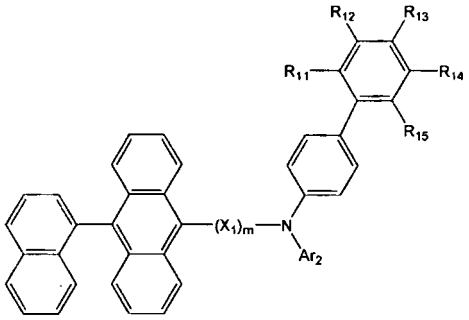
zoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN; a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group; and p and q are each independently 2, 3, or 4.

11. The amine-based compound of claim 1, wherein the amine-based compound is represented by one of Formulae 1A to 1J below:

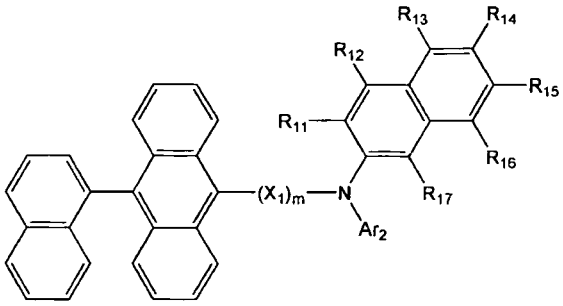
Formula 1A



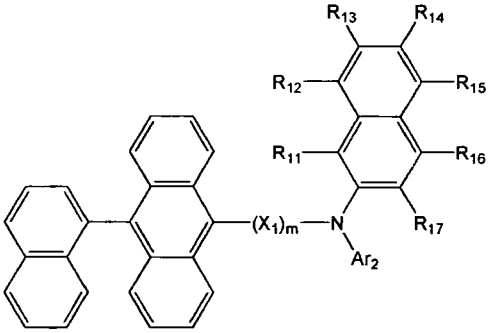
Formula 1B



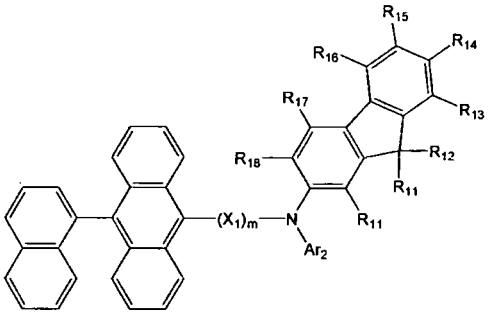
Formula 1C



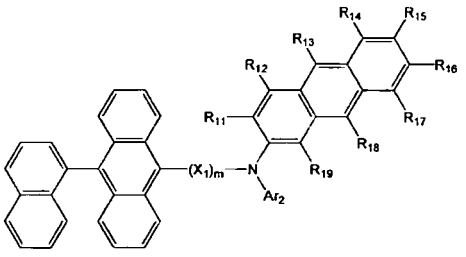
Formula 1D



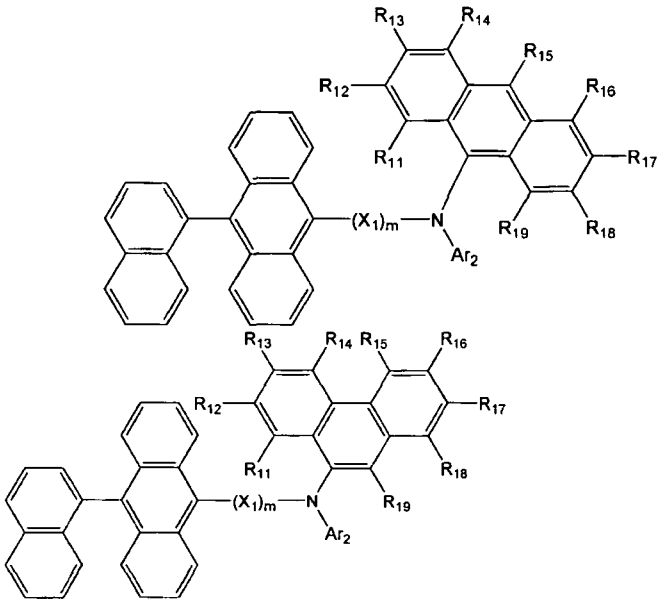
Formula 1E



Formula 1F

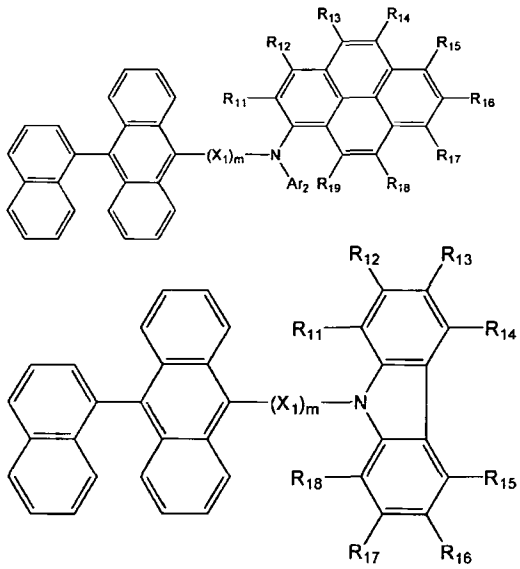


Formula 1G



Formula 1H

Formula 1I



Formula 1J

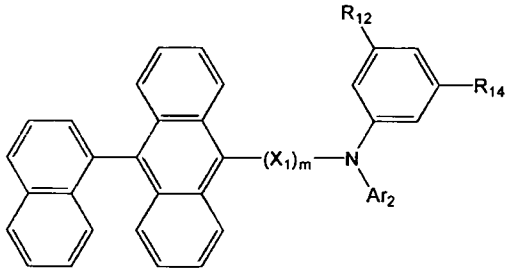
wherein, in Formulae 1A to 1J, Ar_2 is a substituted or unsubstituted phenyl group, a substituted or unsubstituted pentalenyl group, a substituted or unsubstituted indenylene group, a substituted or unsubstituted naphthyl group, a substituted or unsubstituted azulenyl group, a substituted or unsubstituted heptalenyl group, a substituted or unsubstituted indacenyl group, a substituted or unsubstituted acenaphthyl group, a substituted or unsubstituted fluorenyl group, a substituted or unsubstituted phenalenyl group, a substituted or unsubstituted phenanthrenyl group, a substituted or unsubstituted anthryl group, a substituted or unsubstituted fluoranthenyl group, a substituted or unsubstituted triphenylenyl group, a substituted or unsubstituted pyrenyl group, a substituted or unsubstituted chrysenyl group, a substituted or unsubstituted naphthacenyl group, a substituted or unsubstituted picenyl group, a substituted or unsubstituted perylenyl group, a substituted or unsubstituted pentaphenyl group, a substituted or unsubstituted hexacenyl group, a substituted or unsubstituted pyrrolyl group, a substituted or unsubstituted pyrazolyl group, a substituted or unsubstituted imidazolyl group, a substituted or unsubstituted imidazolinyl group, a substituted or unsubstituted imidazopyridinyl group, a substituted or unsubstituted imidazopyrimidinyl group, a substituted or unsubstituted pyridinyl group, a substituted or unsubstituted pyrazinyl group, a substituted or unsubstituted pyrimidinyl group, a substituted or unsubstituted benzoimidazolyl group, a substituted or unsubstituted indolyl group, a

substituted or unsubstituted purinyl group, a substituted or unsubstituted quinolinyl group, a substituted or unsubstituted phthalazinyl group, a substituted or unsubstituted indolizinyl group, a substituted or unsubstituted naphthyridinyl group, a substituted or unsubstituted quinazolinyl group, a substituted or unsubstituted cinolinyl group, a substituted or unsubstituted indazolyl group, a substituted or unsubstituted carbazolyl group, a substituted or unsubstituted phenazinyl group, a substituted or unsubstituted phenanthridinyl group, a substituted or unsubstituted pyranyl group, a substituted or unsubstituted chromenyl group, a substituted or unsubstituted furanyl group, a substituted or unsubstituted benzofuranyl group, a substituted or unsubstituted thiophenyl group, a substituted or unsubstituted benzothiophenyl group, a substituted or unsubstituted isothiazolyl group, a substituted or unsubstituted benzoimidazolyl group, a substituted or unsubstituted isoxazolyl group, a substituted or unsubstituted dibenzothiophenyl group, a substituted or unsubstituted dibenzofuranyl group, a substituted or unsubstituted triazinyl group, a substituted or unsubstituted oxadiazolyl group, a substituted or unsubstituted pyridazinyl group, a substituted or unsubstituted triazolyl group, a substituted or unsubstituted tetrazolyl group, or a substituted or unsubstituted phenanthrolinyl group,

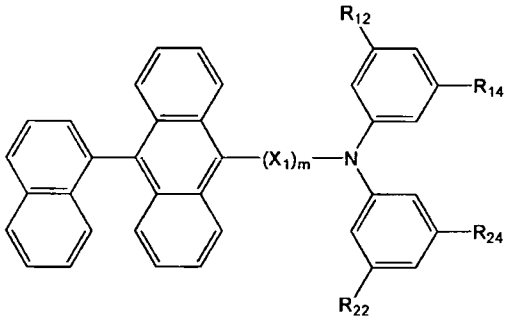
substituents of the substituted phenyl group, the substituted pentalenyl group, the substituted indenyl group, the substituted naphthyl group, the substituted azulenyl group, the substituted heptalenyl group, the substituted indacenyl group, the substituted acenaphthyl group, the substituted fluorenyl group, the substituted phenalenyl group, the substituted phenanthrenyl group, the substituted anthryl group, the substituted fluoranthenyl group, the substituted triphenylenyl group, the substituted pyrenyl group, the substituted chrysenyl group, the substituted naphthacenyl group, the substituted picenyl, the substituted perylenyl group, the substituted pentaphenyl group, the substituted hexacenyl group, the substituted pyrrolyl group, the substituted pyrazolyl group, the substituted imidazolyl group, the substituted imidazolinyl group, the substituted imidazopyridinyl group, the substituted imidazopyrimidinyl group, the substituted pyridinyl group, the substituted pyrazinyl group, the substituted pyrimidinyl group, the substituted benzoimidazolyl group, the substituted indolyl group, the substituted purinyl group, the substituted quinolinyl group, the substituted phthalazinyl group, the substituted indolizinyl group, the substituted naphthyridinyl group, the substituted quinazolinyl group, the substituted cinolinyl group, the substituted indazolyl group, the substituted carbazolyl group, the substituted phenazinyl group, the substituted phenanthridinyl group, the substituted pyranyl group, the substituted chromenyl group, the substituted furanyl group, the substituted benzofuranyl group, the substituted thiophenyl group, the substituted benzothiophenyl group, the substituted isothiazolyl group, the substituted benzoimidazolyl group, the substituted isoxazolyl group, the substituted dibenzothiophenyl group, the substituted dibenzofuranyl group, the substituted triazinyl group, the substituted oxadiazolyl group, the substituted pyridazinyl group, the substituted triazolyl group, the substituted tetrazolyl group, and the substituted phenanthrolinyl group, and R_{11} to R_{19} are each independently a hydrogen atom; a deuterium atom; -F; -Cl; -Br; -I; -CN; a hydroxyl group; $-NO_2$; an amino group; an amidino group; hydrazine; hydrazone; a carboxyl group or a salt thereof; a sulfonic acid group or a salt thereof; phosphoric acid or a salt thereof; a tri(C_6-C_{60}aryl)silyl group; a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, and a C_2-C_{60} alkynyl group; a C_1-C_{60} alkyl group, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group and a C_2-C_{60} alkynyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, and phosphoric acid or a salt thereof; a C_3-C_{60} cycloalkyl group, a C_3-C_{60} cycloalkenyl group, a C_6-C_{60} aryl group, a C_2-C_{60} heteroaryl group, a C_6-C_{60} aralkyl group, a C_6-C_{60} aryloxy group, and a C_6-C_{60} arylthiol group; and a C_3-C_{60} cycloalkyl group, a C_3-C_{60} cycloalkenyl group, a C_6-C_{60} aryl group, a C_2-C_{60} heteroaryl group, a C_6-C_{60} aralkyl group, a C_6-C_{60} aryloxy group, and a C_6-C_{60} arylthiol group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one -F, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group and a C_2-C_{60} heteroaryl group, wherein at least one of R_{11} to R_{15} in Formulae 1A and 1B, at least one of R_{11} to R_{17} in Formulae 1C and 1D, at least one of R_{11} to R_{18} in Formulae 1E and 1J, and at least one of R_{11} to R_{19} in Formula 1F, 1G, 1H and 1I are each independently an electron withdrawing group selected from the group consisting of -F; -CN; $-NO_2$; a C_1-C_{60} alkyl group substituted with at least one -F; a C_2-C_{60} heteroaryl group; and a C_2-C_{60} heteroaryl group substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{60} alkyl group, a C_1-C_{60} alkyl group substituted with at least one -F, a C_1-C_{60} alkoxy group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, a C_6-C_{60} aryl group, and a C_2-C_{60} heteroaryl group.

12. The amine-based compound of claim 11, wherein the amine-based compound is represented by Formula 1A-(1) or 1A-(2) below:

Formula 1A-(1)



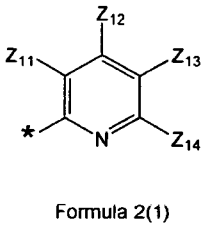
Formula 1A-(2)



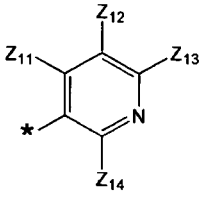
wherein, in Formulae 1A-(1) and 1A-(2), R_{12}, R_{14}, R_{22}, and R_{24} are each independently an electron withdrawing group selected from the group consisting of a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a triazinyl group, a benzoimidazolyl group, and a carbazolyl group; and a pyridinyl group, a pyrazinyl group, a pyrimidinyl group, a quinolinyl group, an isoquinolinyl group, a quinazolinyl group, a phthalazinyl group, a benzoimidazolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN; a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, and a carbazolyl group; and

in Formula 1A-(1) Ar_2 is a substituted or unsubstituted phenyl group, a substituted or unsubstituted biphenyl group, a substituted or unsubstituted naphthyl group, a substituted or unsubstituted anthryl group, a substituted or unsubstituted phenanthrenyl group, a substituted or unsubstituted pyrenyl group, or a substituted or unsubstituted fluorenyl group.

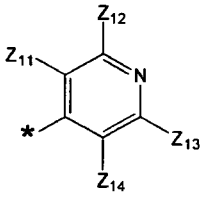
13. The amine-based compound of claim 12, wherein R_{12}, R_{14}, R_{22} and R_{24} are each independently selected from the group consisting of -F; -CN; $-CH_2F$; $-CHF_2$; $-CF_3$; and groups represented by Formulae 2(1) to 2(14) below:



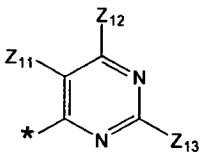
Formula 2(1)



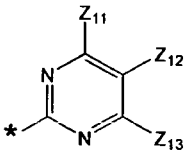
Formula 2(2)



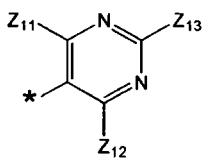
Formula 2(3)



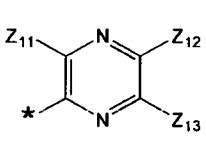
Formula 2(4)



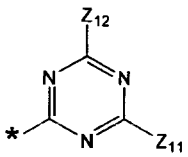
Formula 2(5)



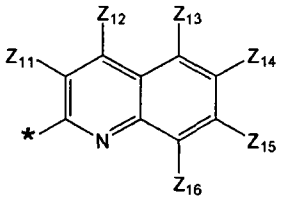
Formula 2(6)



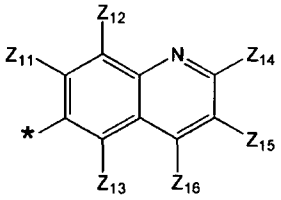
Formula 2(7)



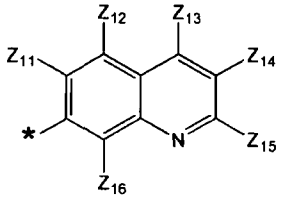
Formula 2(8)



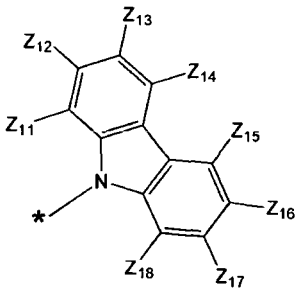
Formula 2(9)



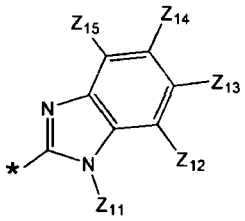
Formula 2(10)



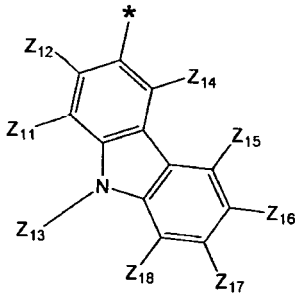
Formula 2(11)



Formula 2(12)



Formula 2(13)

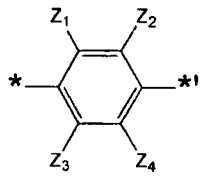


Formula 2(14)

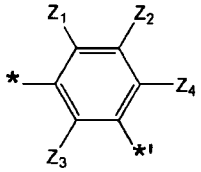
wherein, in Formulae 2(1) to 2(14), Z_{11} to Z_{18} are each independently a hydrogen atom, a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{20}alkyl group, a C_1-C_{20} alkyl group substituted with at least one -F, a C_1-C_{20} alkoxy group, a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a pyridinyl group, a triazinyl group, or a carbazolyl group.

14. The amine-based compound ofany preceding claim, wherein X_1 is a substituted or unsubstituted phenylene group, a substituted or unsubstituted pentalenylene group, a substituted or unsubstituted indenylene group, a substituted or unsubstituted naphthylene group, a substituted or unsubstituted azulenylene group, a substituted or unsubstituted heptalenylene group, a substituted or unsubstituted indacenylene group, a substituted or unsubstituted acenaphthylene group, a substituted or unsubstituted fluorenylene group, a substituted or unsubstituted phenalenylene group, a substituted or unsubstituted phenanthrenylene group, a substituted or unsubstituted anthrylene group, a substituted or unsubstituted fluoranthenylene group, a substituted or unsubstituted triphenylenylene group, a substituted or unsubstituted pyrenylene group, a substituted or unsubstituted chrysenylene group, a substituted or unsubstituted naphthacenylene group, a substituted or unsubstituted picenylene group, a substituted or unsubstituted perylenylene group, a substituted or unsubstituted pentaphenylene group, a substituted or unsubstituted hexacenylene group, a substituted or unsubstituted pyrrolylene group, a substituted or unsubstituted pyrazolylene group, a substituted or unsubstituted imidazolylene group, a substituted or unsubstituted imidazolinylene group, a substituted or unsubstituted imidazopyridinylene group, a substituted or unsubstituted imidazopyrimidinylene group, a substituted or unsubstituted pyridinylene group, a substituted or unsubstituted pyrazinylene group, a substituted or unsubstituted pyrimidinylene group, a substituted or unsubstituted indolylene group, a substituted or unsubstituted purinylene group, a substituted or unsubstituted quinolinylene group, a substituted or unsubstituted phthalazinylene group, a substituted or unsubstituted indolizinylene group, a substituted or unsubstituted naphthyridinylene group, a substituted or unsubstituted quinazolinylene group, a substituted or unsubstituted cinnolinylene group, a substituted or unsubstituted indazolylene group, a substituted or unsubstituted carbazolylene group, a substituted or unsubstituted phenazinylene group, a substituted or unsubstituted phenanthridinylene group, a substituted or unsubstituted pyranylene group, a substituted or unsubstituted chromenylene group, a substituted or unsubstituted furanylene group, a substituted or unsubstituted benzofuranylene group, a substituted or unsubstituted thiophenylene group, a substituted or unsubstituted benzothiophenylene group, a substituted or unsubstituted isothiazolylene group, a substituted or unsubstituted benzoimidazolylene group, a substituted or unsubstituted isoxazolylene group, a substituted or unsubstituted dibenzothiophenylene group, a substituted or unsubstituted dibenzofuranylene group, a substituted or unsubstituted triazinylene group, a substituted or unsubstituted oxadiazolylene group, a substituted or unsubstituted pyridazinylene group, a substituted or unsubstituted triazolylene group, or a substituted or unsubstituted tetrazolylene group.

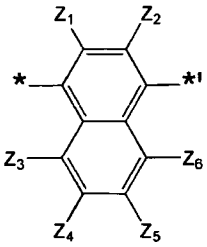
15. The amine-based compound of claim 14, wherein X_1 is a group represented by one of Formulae 5(1) to 5(16) below:



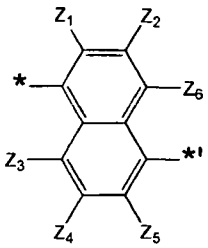
Formula 5(1)



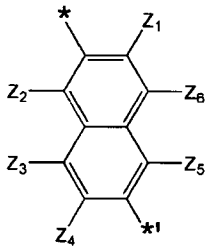
Formula 5(2)



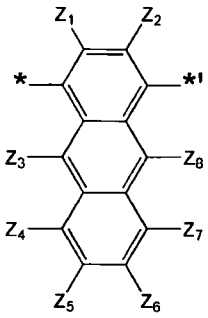
Formula 5(3)



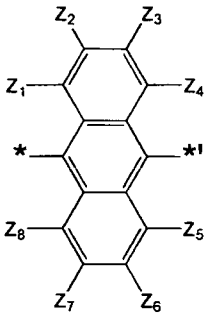
Formula 5(4)



Formula 5(5)



Formula 5(6)



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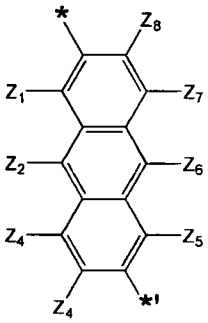

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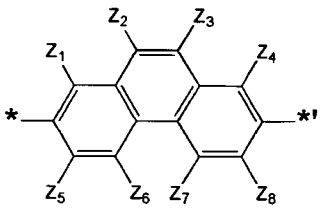
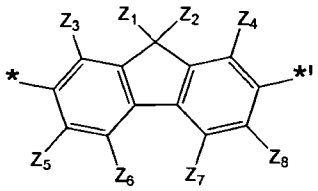
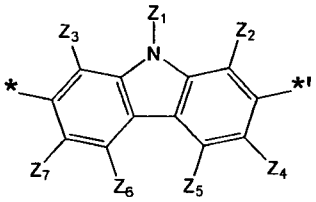


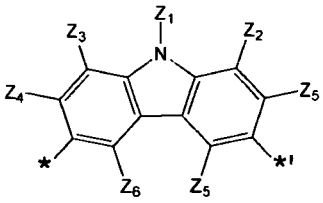
Formula 5(9)



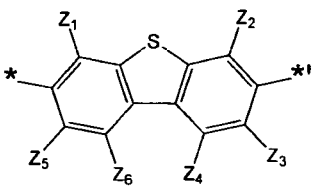
Formula 5(10)



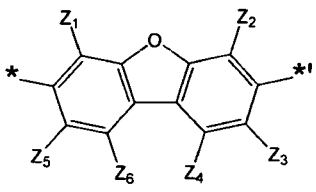
Formula 5(11)



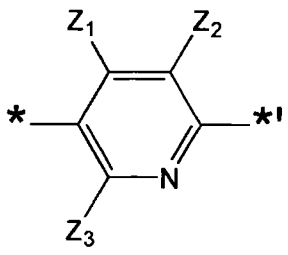
Formula 5(12)



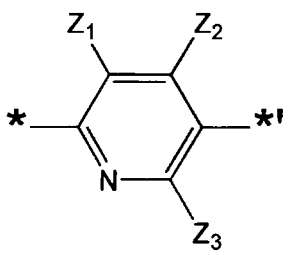
Formula 5(13)



Formula 5(14)



Formula 5(15)



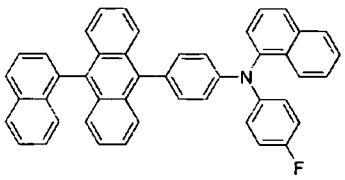
Formula 5(16)

wherein, in Formulae 5(1) to 5(16), Z_1 to Z_8 are each independently one of a hydrogen atom; a deuterium atom; -F; -Cl; -Br; -I; -CN; a hydroxyl group; $-NO_2$; an amino group; an amidino group; hydrazine; hydrazone; a carboxyl group or a salt thereof; a sulfonic acid group or a salt thereof; phosphoric acid or a salt thereof; a C_1-C_{20} alkyl group; a C_1-C_{20} alkoxy group; a C_1-C_{20} alkyl group and a C_1-C_{20} alkoxy groups that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, sulfonic acid group or a salt thereof, and phosphoric acid or a salt thereof; a C_6-C_{20} aryl group; a C_2-C_{20} heteroaryl group; and a C_6-C_{20} aryl group and a C_2-C_{20} heteroaryl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, phosphoric acid or a salt thereof, a C_1-C_{20} alkyl group, a C_1-C_{20} alkoxy group, a C_6-C_{20} aryl group and a C_2-C_{20} heteroaryl group;
wherein indicates a binding site to anthracene in Formula 1; and
wherein *' indicates a binding site to N in Formula 1.

16. The amine-based compound of claim 15, wherein Z_1 to Z_8 are each independently one of a hydrogen atom; a deuterium atom; -F; -Cl; -Br; -I; -CN; a hydroxyl group; $-NO_2$; an amino group; an amidino group; hydrazine; hydrazone; a carboxyl group or a salt thereof; a sulfonic acid group or a salt thereof; a phosphoric acid or a salt thereof; a methyl group, an ethyl group, a propyl group, a butyl group, and a pentyl group; a methoxy group, an ethoxy group, a propoxy group, a butoxy group, and a pentoxy group; a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, a methoxy group, an ethoxy group, a propoxy group, a butoxy group, and a pentoxy group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, and a phosphoric acid or a salt thereof; a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, and a fluorenyl group; a pyridinyl group, a pyrimidinyl group, a triazinyl group, a quinolyl group, and a carbazolyl group; a phenyl group, a naphthyl group, an anthryl group, a phenanthrenyl group, a pyrenyl group, a fluorenyl group, a pyridinyl group, a pyrimidinyl group, a triazinyl group, a quinolyl group, and a carbazolyl group that are substituted with at least one of a deuterium atom, -F, -Cl, -Br, -I, -CN, a hydroxyl group, $-NO_2$, an amino group, an amidino group, hydrazine, hydrazone, a carboxyl group or a salt thereof, a sulfonic acid group or a salt thereof, a phosphoric acid or a salt thereof, a C_1-C_{20}alkyl group, and a C_1-C_{20}alkoxy group.

17. The amine-based compound of any preceding claim, wherein m is 1, 2, or 3.

18. The amine-based compound of any preceding claim, wherein the amine-based compound is one of Compounds 1 to 109 below:



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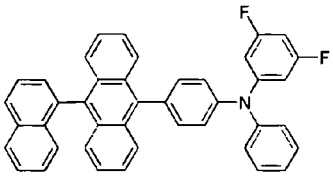

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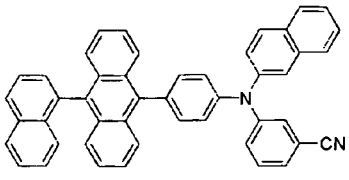

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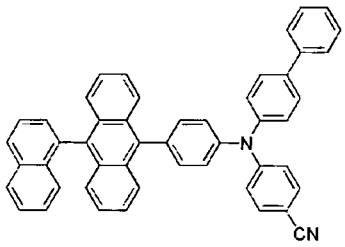

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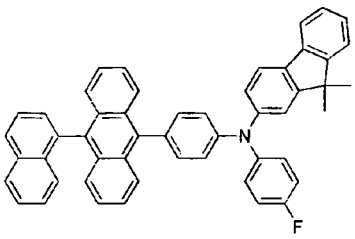

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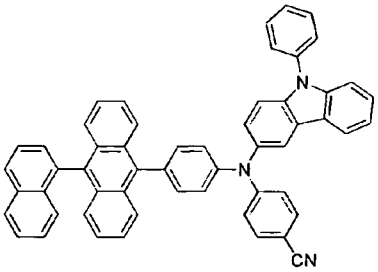

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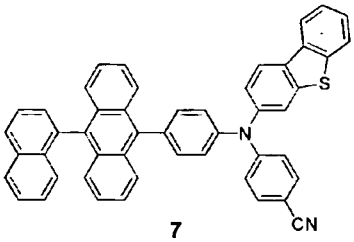

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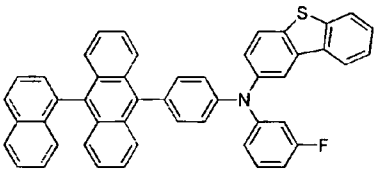

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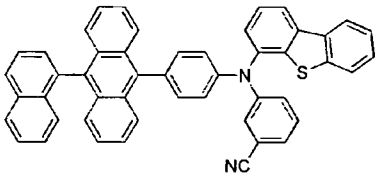

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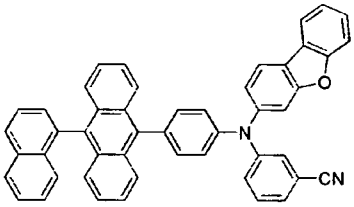

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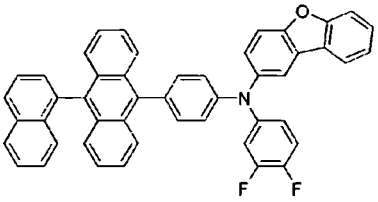

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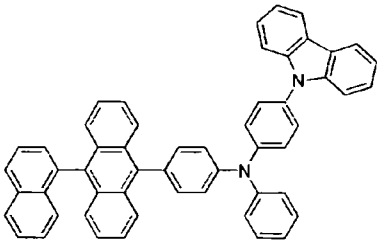

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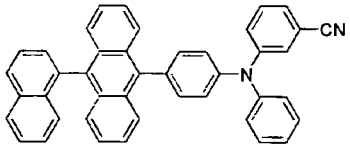

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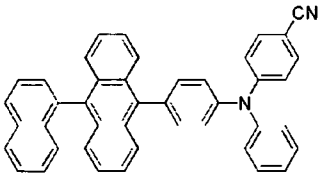

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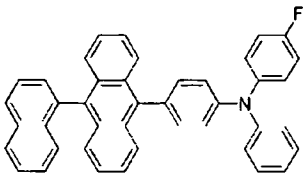

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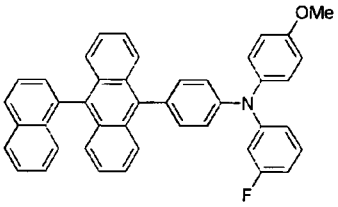

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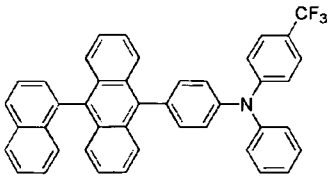

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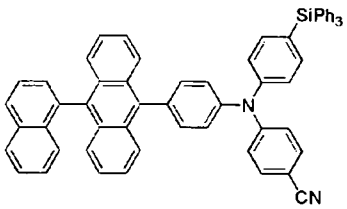

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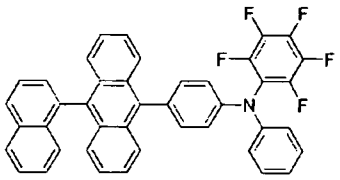

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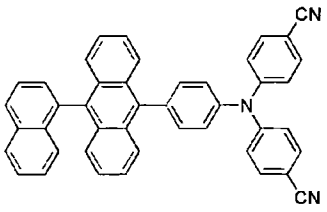

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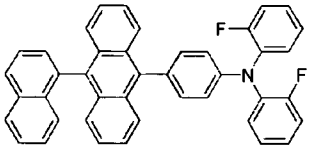

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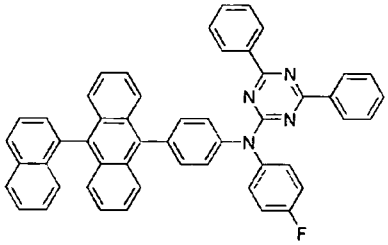

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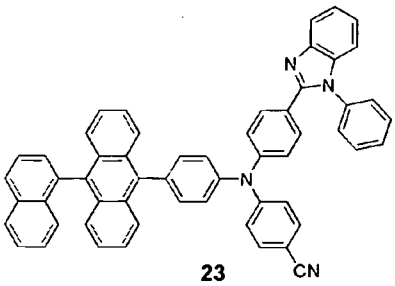

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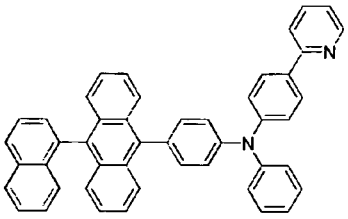

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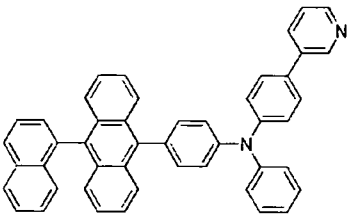

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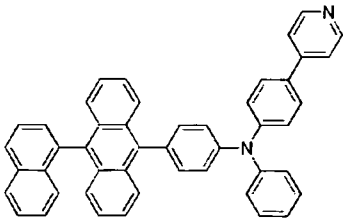

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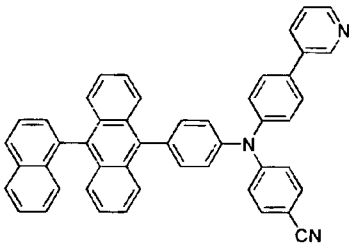

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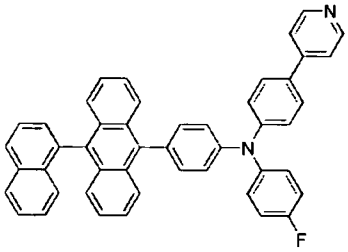

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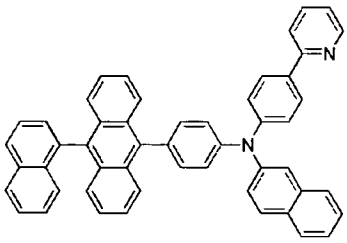

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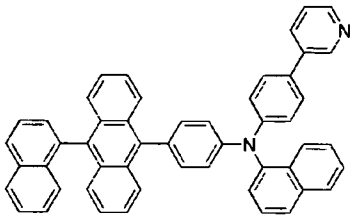

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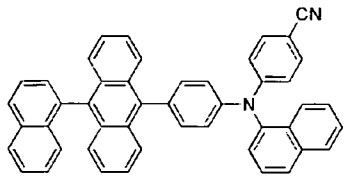

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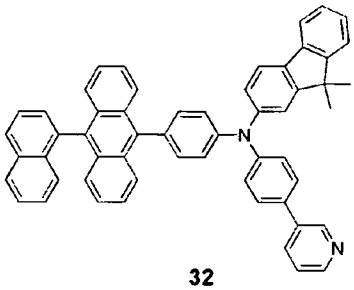

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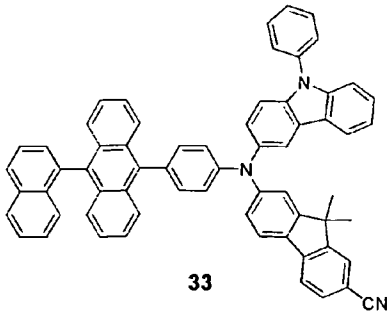

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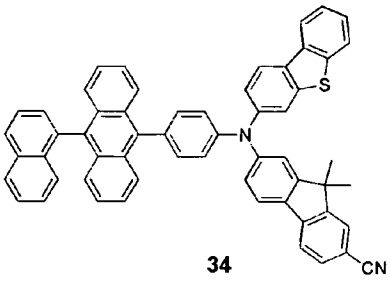

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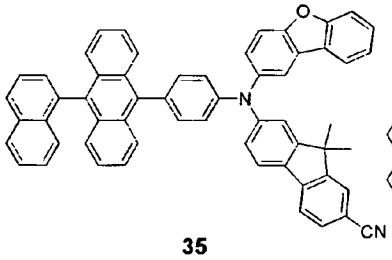

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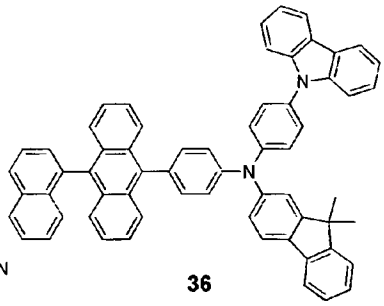

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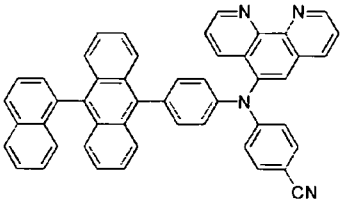

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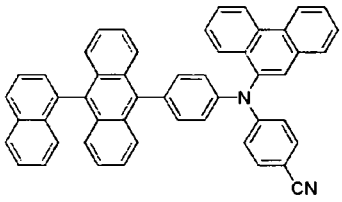

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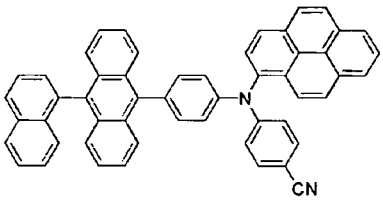

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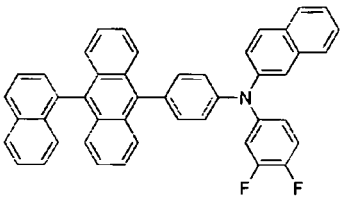

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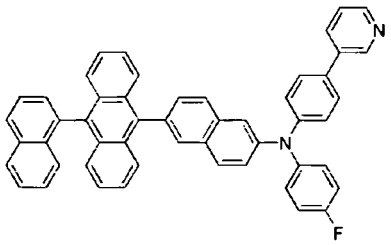

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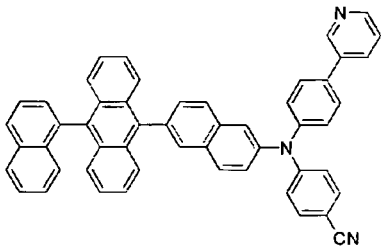

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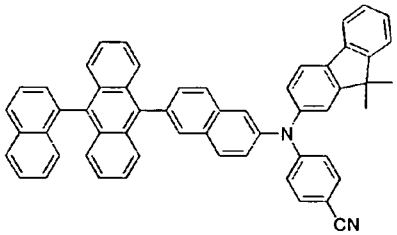

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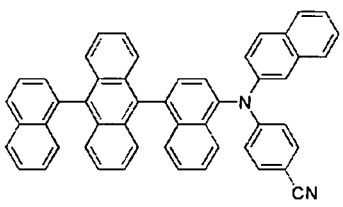

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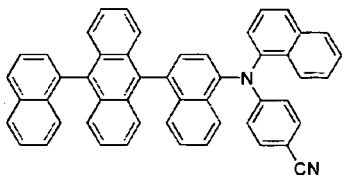

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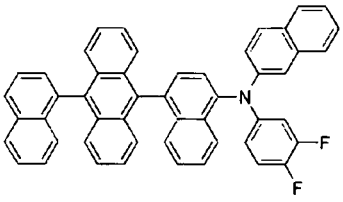

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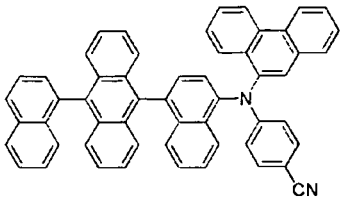

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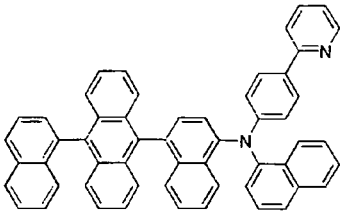

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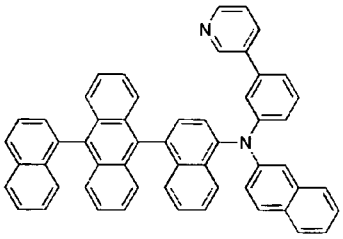

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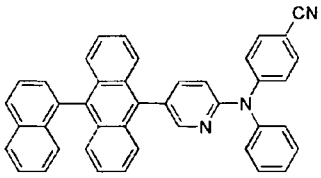

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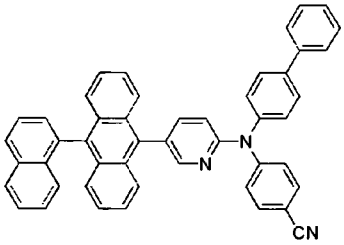

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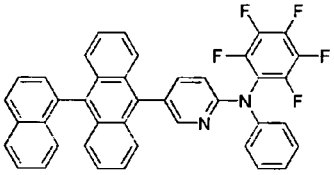

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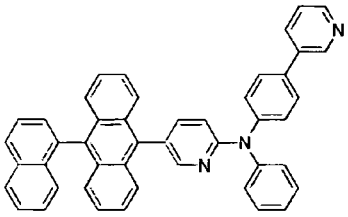

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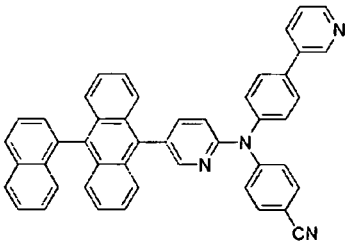

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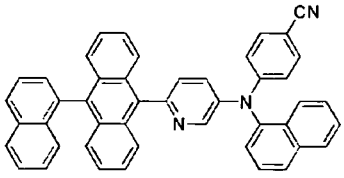

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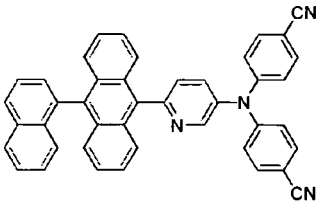

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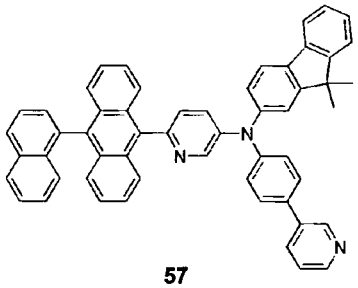

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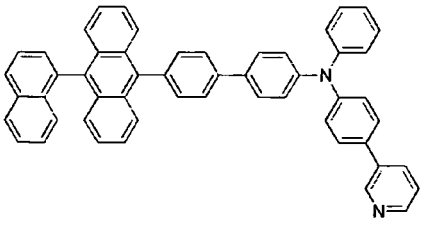

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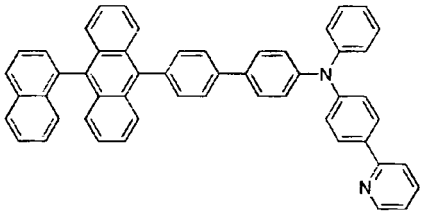

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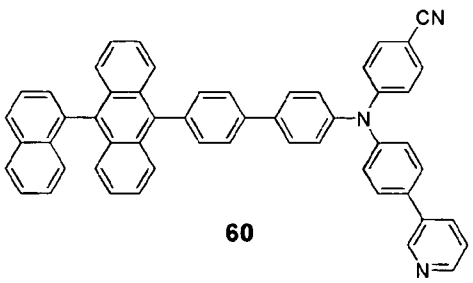

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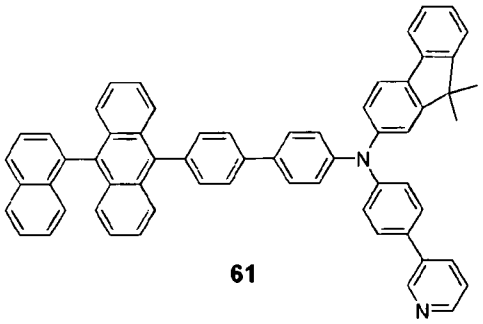

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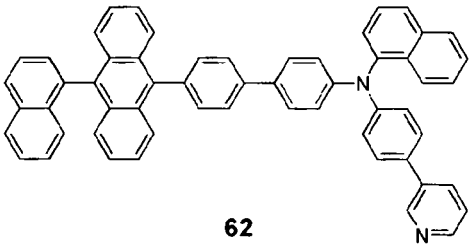

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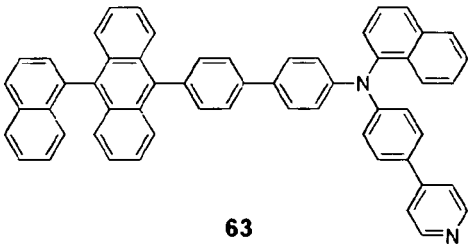

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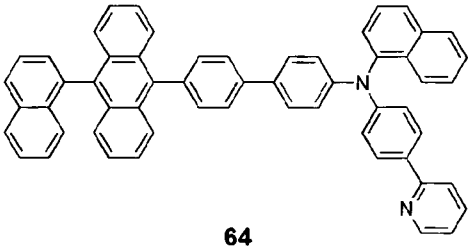

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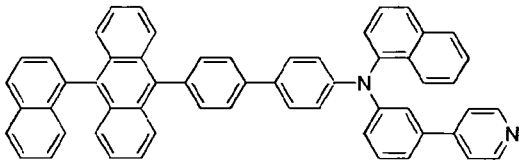

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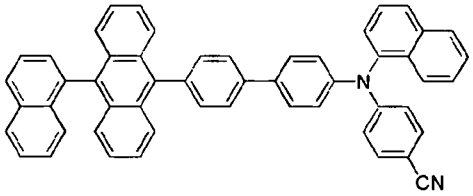

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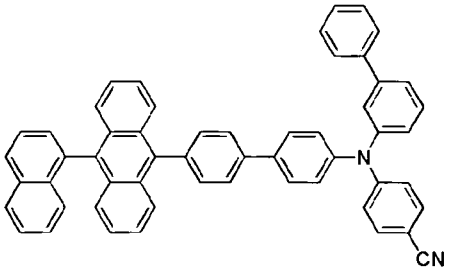

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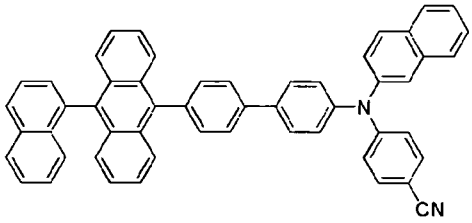

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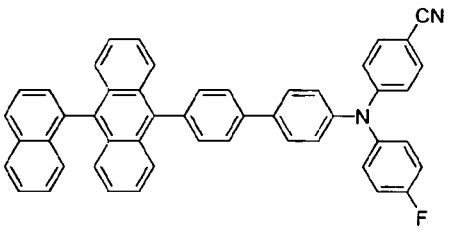

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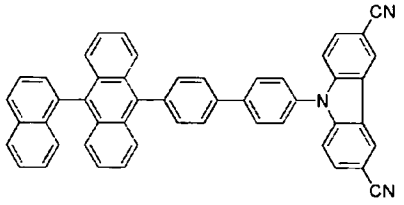

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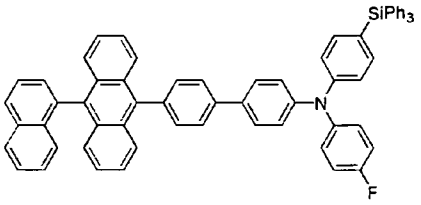

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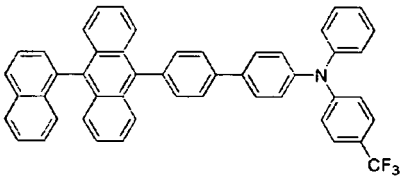

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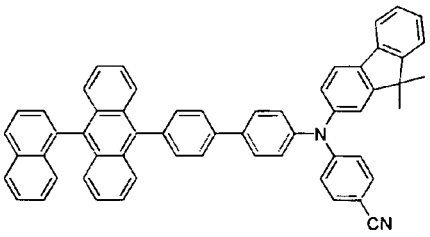

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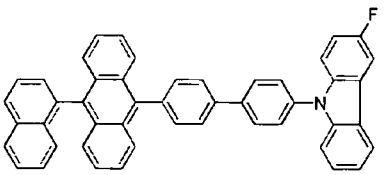

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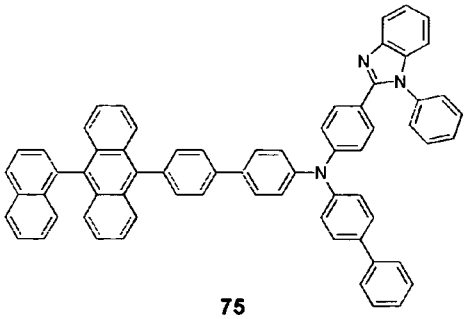

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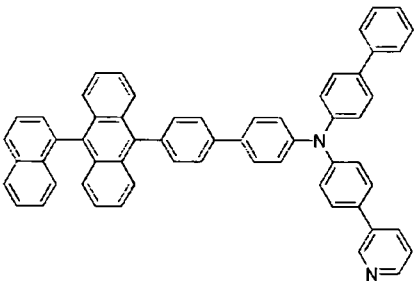

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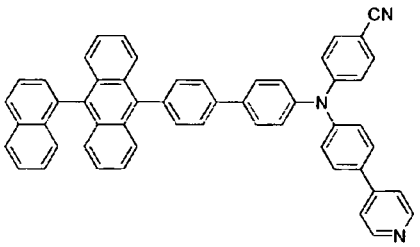

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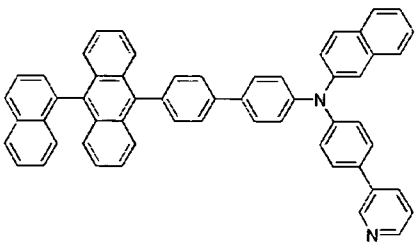

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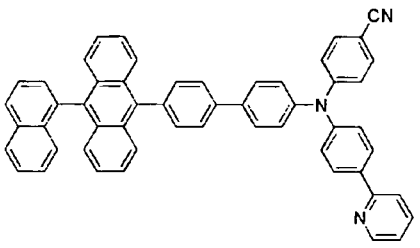

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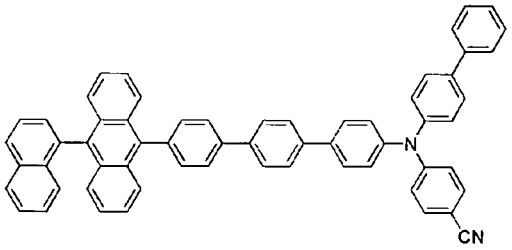

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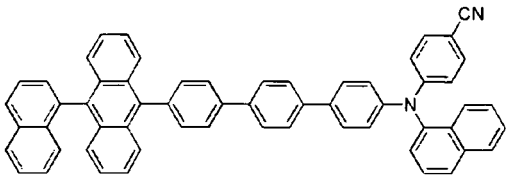

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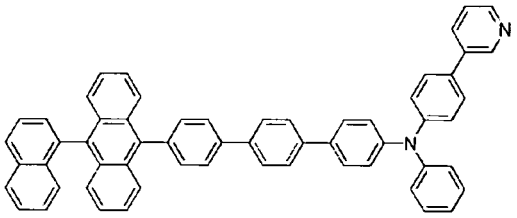

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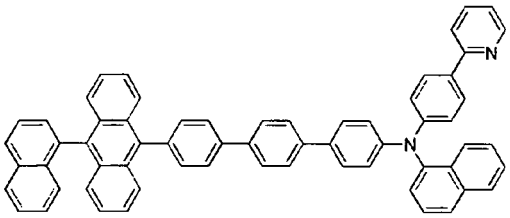

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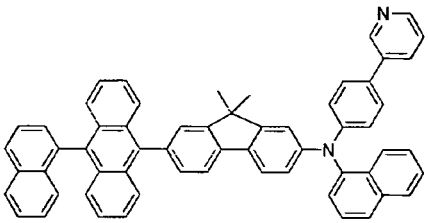

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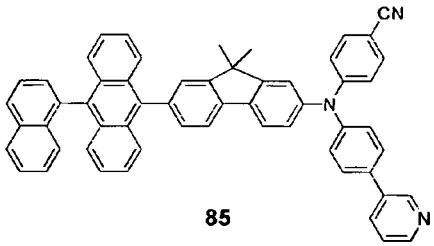

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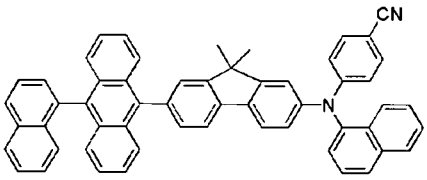

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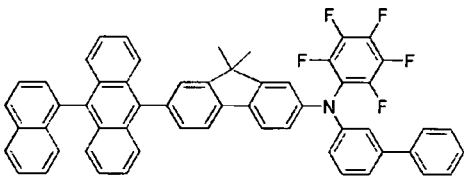

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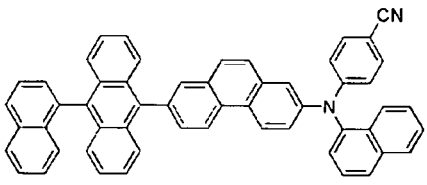

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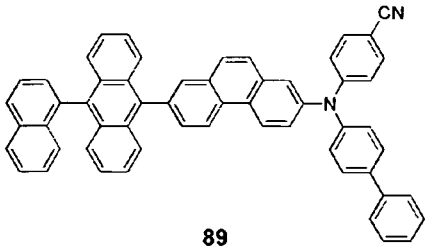

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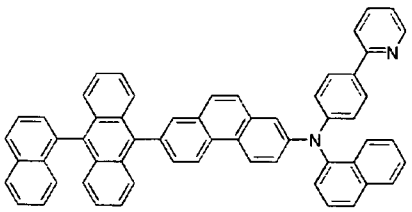

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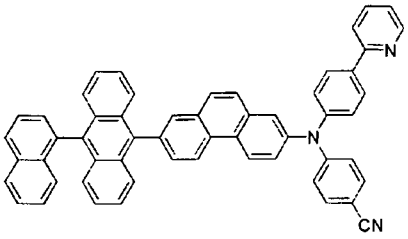

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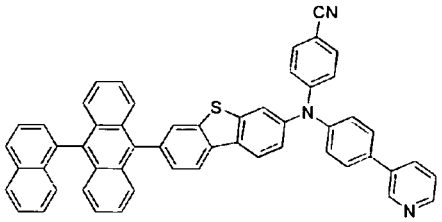

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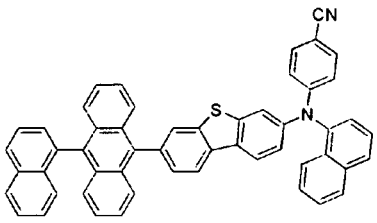

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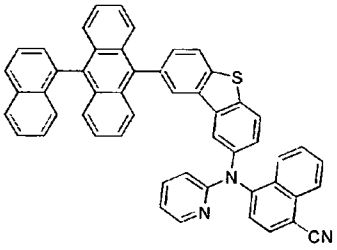

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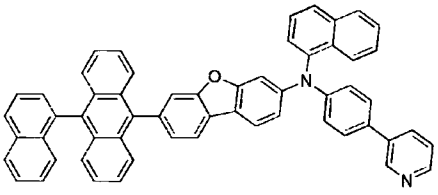

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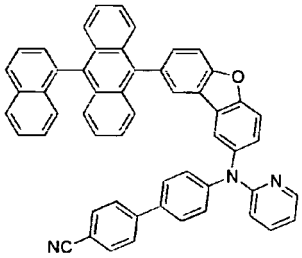

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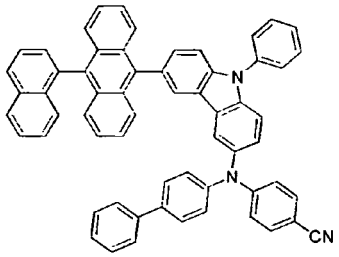

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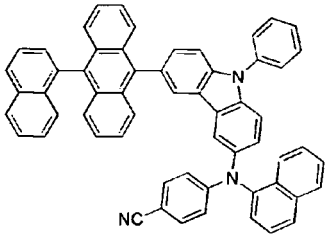

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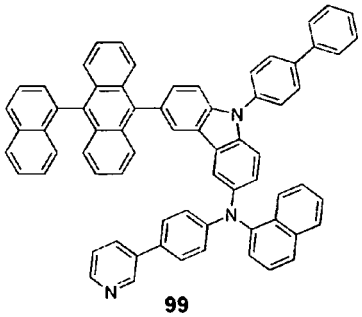

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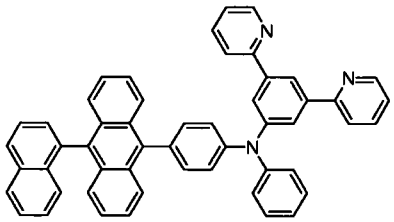

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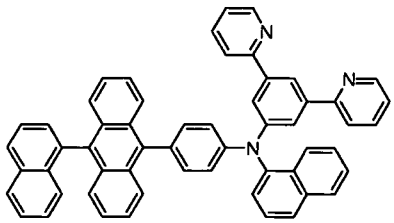

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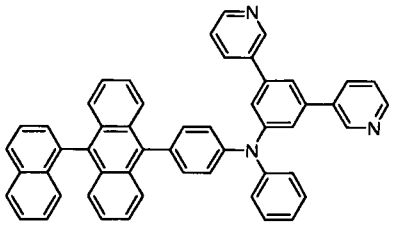

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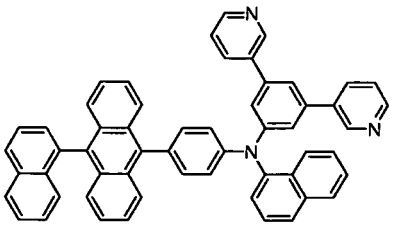

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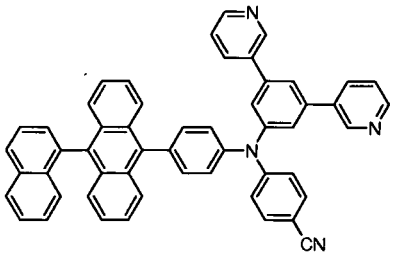

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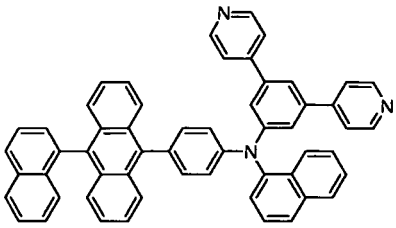

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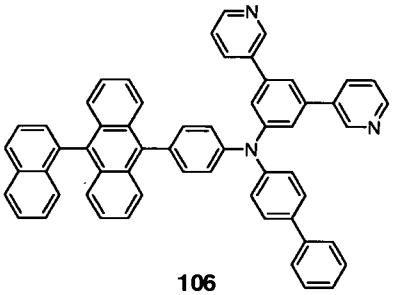

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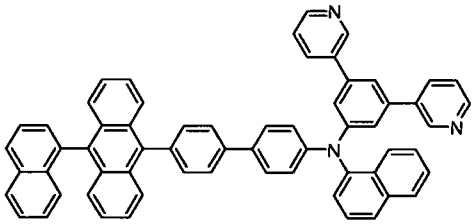

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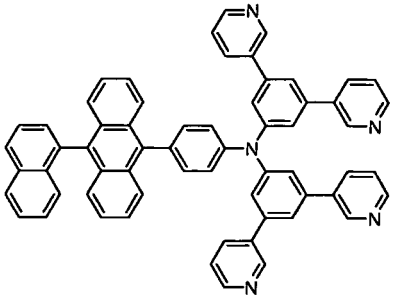

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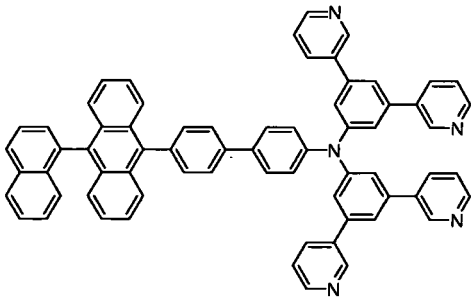

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19. An organic light-emitting diode comprising a first electrode, a second electrode disposed opposite to the first electrode, and an organic layer disposed between the first electrode and the second electrode, the organic layer comprising at least one of the amine-based compounds of claims 1 to 18.

20. The organic light-emitting diode of claim 19, wherein the organic layer comprises at least one of a hole injection layer, a hole transport layer, a functional layer having both hole injection and hole transport capabilities, a buffer layer, an electro blocking layer, an emission layer, a hole blocking layer, an electron transport layer, an electron injection layer, and a functional layer having both electron injection and electron transport capabilities.

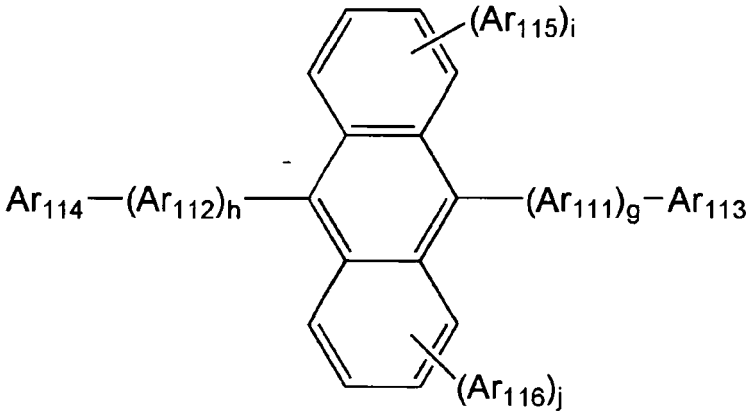
21. The organic light-emitting diode of claim 20, wherein the organic layer comprises an electron transport layer, and the amine-based compound is included in the electron transport layer, preferably
wherein the electron transport layer further comprises a metal complex.

22. The organic light-emitting diode of claim 20, wherein the organic layer comprises an emission layer, and the amine-

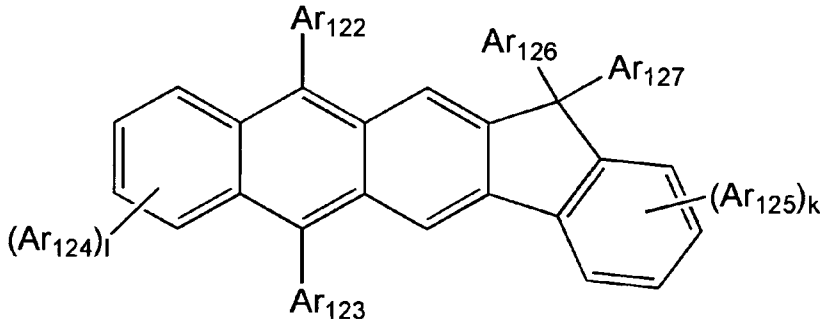
based compound is included in the emission layer, preferably.
wherein the amine-based compound in the emission layer serves as a host, and the emission layer further comprises a blue fluorescent dopant.

23. The organic light-emitting diode of claim 22, wherein the amine-based compound in the emission layer serves as a dopant, and the emission layer further comprises at least one of an anthracene-based compound represented by Formula 400 below and an anthracene-based compound represented by Formula 401 below:

Formula 400



Formula 401



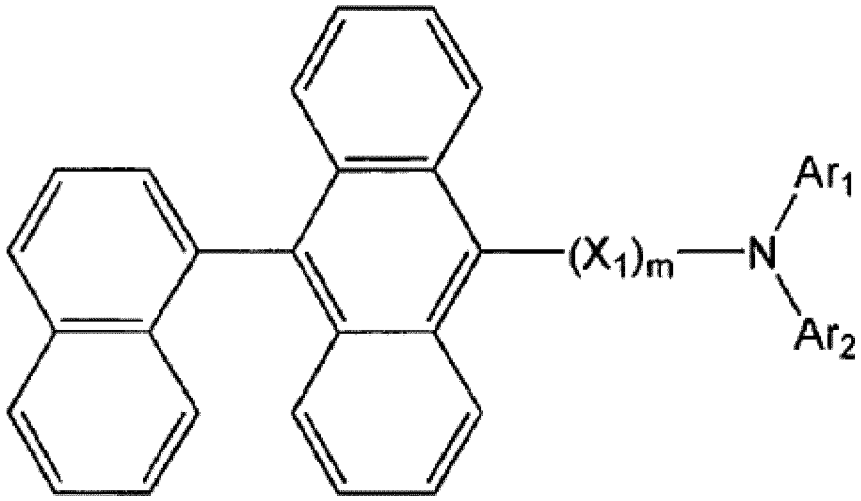
wherein, in Formulae 400 and 401, Arm and Ar_{112} are each independently a substituted or unsubstituted C_6-C_{60} arylene group; Ar_{113} to Ar_{116}, and Ar_{122} to Ar_{125} are each independently a substituted or unsubstituted C_1-C_{10} alkyl group, or a substituted or unsubstituted C_6-C_{60} aryl group; Ar_{126} and Ar_{127} are each independently a C_1-C_{10} alkyl group; and g, h, i, j, k, and 1 are each independently an integer from 0 to 4.

24. The organic light-emitting diode of claim 20, wherein the organic layer comprises at least one of a hole injection layer, a hole transport layer, and a functional layer having both hole injection and hole transport capabilities, and the at least one of the hole injection layer, the hole transport layer, and the functional layer having both hole injection and hole transport capabilities comprises a p-dopant.

Patentansprüche

1. Aminbasierte Verbindung, wiedergegeben durch nachstehende Formel 1 :

Formel 1



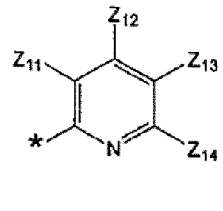
wobei, in Formel 1, Ar_1 und Ar_2 jeweils unabhängig voneinander ein substituierter oder unsubstituierter C_6-C_{60}-Arylrest oder ein substituierter oder unsubstituierter C_2-C_{60}-Heteroarylrest sind; und wobei Ar_1 und Ar_2 gegebenenfalls über eine Einfachbindung verbunden sind;

X_1 ein substituierter oder unsubstituierter C_6-C_{60}-Arylenrest oder ein substituierter oder unsubstituierter C_2-C_{60}-Heteroarylenrest ist;
m eine ganze Zahl von 1 bis 5 ist; und
mindestens ein Substituent von jedem des substituierten C_6-C_{60}-Arylrestes, des substituierten C_2-C_{60}-Heteroarylrestes, des substituierten C_6-C_{60}-Arylenrestes und des substituierten C_2-C_{60}-Heteroarylenrestes eines ist von einem Deuteriumatom; -F; -Cl; -Br; -I; -CN; einer Hydroxylgruppe; $-NO_2$; einer Aminogruppe; einer Amidinogruppe; Hydrazin; Hydrazon; einer Carboxylgruppe oder einem Salz davon; einer Sulfonsäuregruppe oder einem Salz davon; einer Phosphorsäure oder einem Salz davon; einem Tri(C_6-C_{60}-aryl)silylrest; einem C_1-C_{60}-Alkylrest, einem C_1-C_{60}-Alkoxyrest, einem C_2-C_{60}-Alkenylrest, einem C_2-C_{60}-Alkinylrest; einem C_1-C_{60}-Alkylrest, einem C_1-C_{60}-Alkoxyrest, einem C_2-C_{60}-Alkenylrest und einem C_2-C_{60}-Alkinylrest, der substituiert ist mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, einer Phosphorsäure oder einem Salz davon; einem C_3-C_{60}-Cycloalkylrest, einem C_3-C_{60}-Cycloalkenylrest, einem C_6-C_{60}-Arylrest, einem C_2-C_{60}-Heteroarylrest, einem C_6-C_{60}-Aralkylrest, einem C_6-C_{60}-Aryloxyrest, einem C_6-C_{60}-Arylthiolrest; einem C_3-C_{60}-Cycloalkylrest, einem C_3-C_{60}-Cycloalkenylrest, einem C_6-C_6-Arylrest, einem C_2-C_{60}-Heteroarylrest, einem C_6-C_{60}-Aralkylrest, einem C_6-C_{60}-Aryloxyrest und einem C_6-C_{60}-Arylthiolrest, der substituiert ist mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, einer Phosphorsäure oder einem Salz davon, einem C_1-C_{60}-Alkylrest, einem mit mindestens einem Fluoratom (F) substituierten C_1-C_{60}-Alkylrest, einem C_1-C_{60}-Alkoxyrest; einem C_2-C_{60}-Alkenylrest, einem C_2-C_{60}-Alkinylrest; einem C_6-C_{60}-Arylrest und einem C_2-C_{60}-Heteroarylrest,
wobei mindestens eines von Ar_1 und Ar_2 ein C_6-C_{60}-Arylrest ist, substituiert mit mindestens einer elektronenziehenden Gruppe, ausgewählt aus -F; -CN; $-NO_2$; einem mit mindestens einem-F substituierten C_1-C_{60}-Alkylrest; einem C_2-C_{60}-Heteroarylrest; und einem C_2-C_{60}-Heteroarylrest, substituiert mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, einer Phosphorsäure oder einem Salz davon, einem C_1-C_{60}-Alkylrest, einem mit mindestens einem -F substituierten C_1-C_{60}-Alkylrest, einem C_1-C_{60}-Alkoxyrest, einem C_2-C_{60}-Alkenylrest, einem C_2-C_{60}-Alkinylrest, einem C_6-C_{60}-Arylrest und einem C_2-C_{60}-Heteroarylrest.

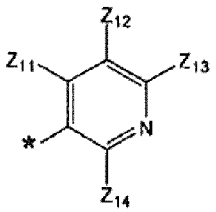
2. Aminbasierte Verbindung nach Anspruch 1, wobei mindestens eine elektronenziehende Gruppe ausgewählt ist aus: -F; -CN; $-NO_2$, einem mit mindestens einem -F substituierten C_1-C_{20}-Alkylrest; einem C_2-C_{20}-Heteroarylrest, umfassend einen ein N-Atom enthaltenden Ring; und einem C_2-C_{20}-Heteroarylrest, der einen ein N-Atom enthaltenden

Ring umfasst und substituiert ist mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, Phosphorsäure oder einem Salz davon, einem C_1-C_{20}-Alkylrest, einem mit mindestens einem -F substituierten C_1-C_{20}-Alkylrest, einem C_1-C_{20}-Alkoxyrest, einem C_6-C_{20}-Arylrest und einem C_2-C_{20}-Heteroarylrest.

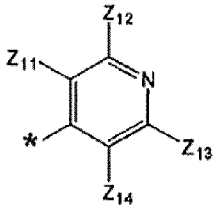
3. Aminbasierte Verbindung nach Anspruch 1 oder Anspruch 2, wobei mindestens eine elektronenziehende Gruppe ausgewählt ist aus -F; -CN; $-CH_2F$; $-CHF_2$; $-CF_3$ und den durch die nachstehenden Formeln 2(1) bis 2(14) dargestellten Resten:



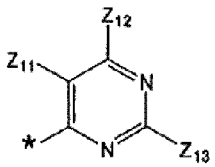
Formel 2(1)



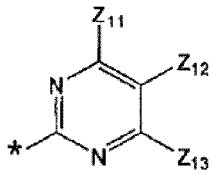
Formel 2(2)



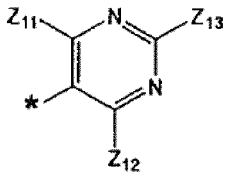
Formel 2(3)



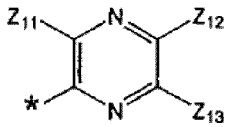
Formel 2(4)



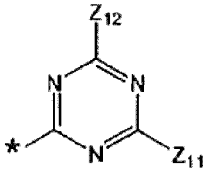
Formel 2(5)



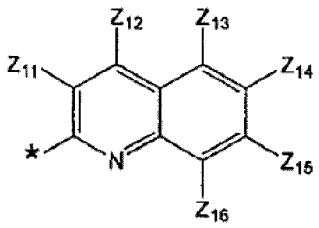
Formel 2(6)



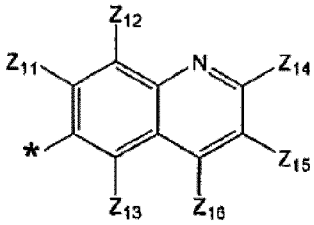
Formel 2(7)



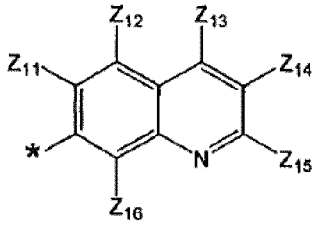
Formel 2(8)



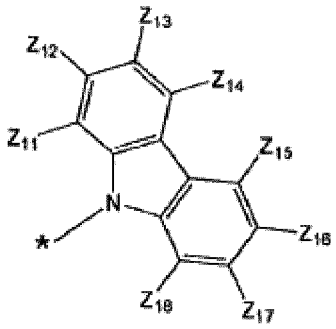
Formel 2(9)



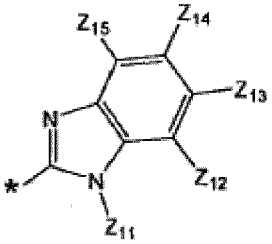
Formel 2(10)



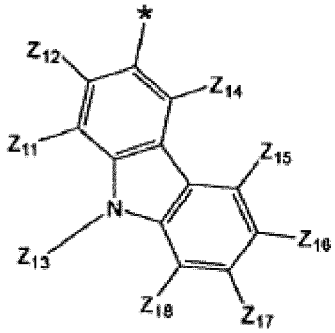
Formel 2(11)



Formel 2(12)



Formel 2(13)



Formel 2(14)

wobei, in den Formeln 2(1) bis 2(14), Z_{11} bis Z_{18} jeweils unabhängig voneinander ein Wasserstoffatom, ein Deuteriumatom, -F, -Cl, -Br, -I, -CN, eine Hydroxylgruppe, $-NO_2$, eine Aminogruppe, eine Amidinogruppe, Hydrazin, Hydrazon, eine Carboxylgruppe oder ein Salz davon, eine Sulfonsäuregruppe oder ein Salz davon, Phosphorsäure

oder ein Salz davon, ein C_1-C_{20}-Alkylrest, ein mit mindestens einem -F substituierter C_1-C_{20}-Alkylrest, ein C_1-C_{20}-Alkoxyrest, eine Phenylgruppe, eine Naphthylgruppe, eine Anthrylgruppe, eine Phenanthrenylgruppe, eine Pyrenylgruppe, eine Pyridinylgruppe, eine Triazinylgruppe oder eine Carbazolylgruppe sind.

4. Aminbasierte Verbindung nach einem vorangehenden Anspruch, wobei Ar_1 und Ar_2 jeweils unabhängig voneinander ein C_6-C_{60}-Arylrest sind, ausgewählt aus einer substituierten oder unsubstituierten Phenylgruppe, einer substituierten oder unsubstituierten Naphthylgruppe, einer substituierten oder unsubstituierten Fluorenylgruppe, einer substituierten oder unsubstituierten Phenanthrenylgruppe, einer substituierten oder unsubstituierten Anthrylgruppe, einer substituierten oder unsubstituierten Triphenylenylgruppe, einer substituierten oder unsubstituierten Pyrenylgruppe, einer substituierten oder unsubstituierten Chrysenylgruppe, einer substituierten oder unsubstituierten Pyridinylgruppe, einer substituierten oder unsubstituierten Pyrazinylgruppe, einer substituierten oder unsubstituierten Pyrimidinylgruppe, einer substituierten oder unsubstituierten Chinolylgruppe, einer substituierten oder unsubstituierten Carbazolylgruppe, einer substituierten oder unsubstituierten Triazinylgruppe, einer substituierten oder unsubstituierten Dibenzothiophenylgruppe, einer substituierten oder unsubstituierten Dibenzofuranylgruppe oder einer substituierten oder unsubstituierten Phenanthrolinylgruppe.

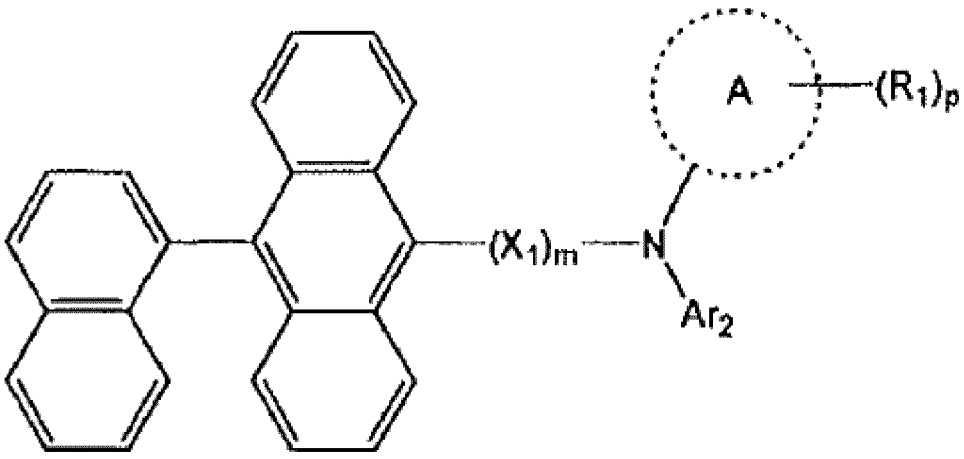
5. Aminbasierte Verbindung nach einem vorangehenden Anspruch, wobei Ar_1 und Ar_2 über eine Einfachbindung verbunden sind.

6. Aminbasierte Verbindung nach einem vorangehenden Anspruch, wobei mindestens eines von Ar_1 und Ar_2 ein mit mindestens zwei elektronenziehenden Gruppen substituierter C_6-C_{60}-Arylrest ist.

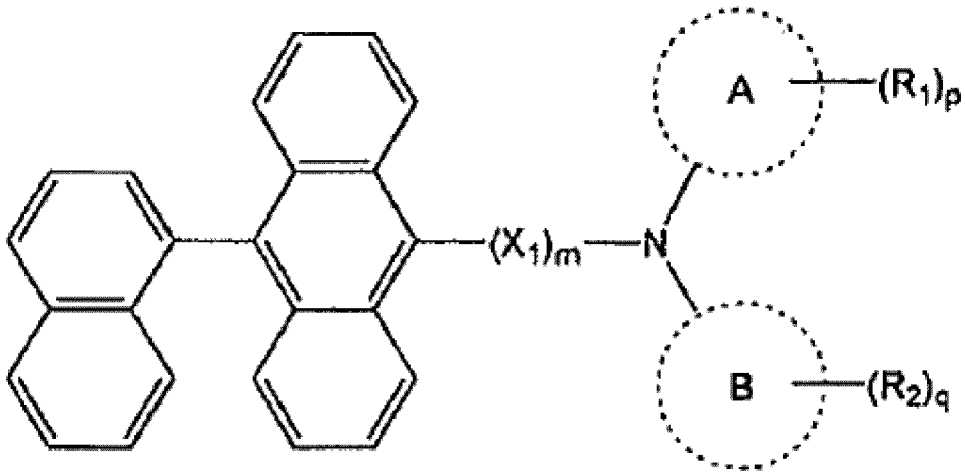
7. Aminbasierte Verbindung nach einem vorangehenden Anspruch, wobei mindestens eines von Ar_1 und Ar_2 eine Phenylgruppe, eine Biphenylgruppe, eine Naphthylgruppe, eine Anthrylgruppe, eine Phenanthrenylgruppe, eine Pyrenylgruppe oder eine Fluorenylgruppe ist, die mit mindestens zwei elektronenziehenden Gruppen substituiert ist; und
die elektronenziehenden Gruppen jeweils unabhängig voneinander ausgewählt sind aus einer Pyridinylgruppe, einer Pyrazinylgruppe, einer Pyrimidinylgruppe, einer Chinolinylgruppe, einer Isochinolinylgruppe, einer Chinazolinylgruppe, einer Triazinylgruppe, einer Benzimidazolylgruppe und einer Carbazolylgruppe; und einer Pyridinylgruppe, einer Pyrazinylgruppe, einer Pyrimidinylgruppe, einer Chinolinylgruppe, einer Isochinolinylgruppe, einer Chinazolinylgruppe, einer Phthalazinylgruppe, einer Benzimidazolylgruppe und einer Carbazolylgruppe, die substituiert sind mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN; einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, einer Phosphorsäure oder einem Salz davon, einem C_1-C_{20}-Alkylrest, einem mit mindestens einem -F substituierten C_1-C_{20}-Alkylrest, einem C_1-C_{20}-Alkoxyrest, einer Phenylgruppe, einer Naphthylgruppe, einer Anthrylgruppe, einer Phenanthrenylgruppe, einer Pyrenylgruppe, einer Pyridinylgruppe, einer Triazinylgruppe und einer Carbazolylgruppe.

8. Aminbasierte Verbindung nach Anspruch 1, wobei die aminbasierte Verbindung durch die nachstehende Formel 1(1) oder 1 (2) wiedergegeben wird:

Formel 1(1)



Formel 1(2)



wobei, in Formel 1(1), Ar_2 ein substituierter oder unsubstituierter C_6-C_{20}-Arylrest oder ein substituierter oder unsubstituierter C_2-C_{20}-Heteroarylrest ist; und
in den Formeln 1(1) und 1(2) die Ringgruppen A und B jeweils unabhängig voneinander ein substituierter C_6-C_{20}-Arylrest sind;
R_1 und R_2 jeweils unabhängig voneinander eine elektronenziehende Gruppe sind, ausgewählt aus -F, -CN, $-NO_2$; einem mit mindestens einem -F substituierten C_1-C_{60}-Alkylrest; einem C_2-C_{60}-Heteroarylrest und einem C_2-C_{60}-Heteroarylrest, substituiert mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, einer Phosphorsäure oder einem Salz davon, einem C_1-C_{60}-Alkylrest, einem mit mindestens einem -F substituierten C_1-C_{60}-Alkylrest, einem C_1-C_{60}-Alkoxyrest, einem C_2-C_{60}-Alkenylrest, einem C_2-C_{60}-Alkinylrest; einem C_6-C_{60}-Arylrest und einem C_2-C_{60}-Heteroarylrest; und
p und q jeweils unabhängig voneinander eine ganze Zahl von 1 bis 9 sind.

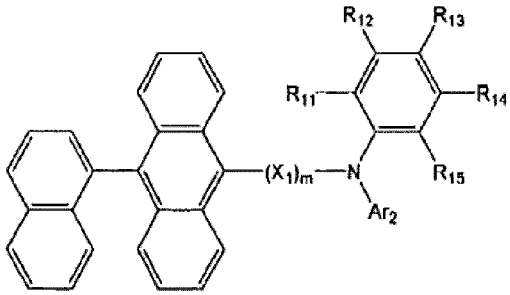
9. Aminbasierte Verbindung nach Anspruch 8, wobei die aminbasierte Verbindung durch die Formel 1 (1) wiedergegeben wird, wobei mindestens eines der R_1 -CN ist; oder durch die Formel 1(2) wiedergegeben wird, wobei mindestens eines der R_1 und eines der R_2 -CN ist.

10. Aminbasierte Verbindung nach Anspruch 7, wobei die aminbasierte Verbindung durch die Formel 1(1) wiedergegeben wird, wobei die Ringgruppe A eine substituierte Phenylgruppe, eine substituierte Biphenylgruppe, eine substituierte Naphthylgruppe, eine substituierte Anthrylgruppe, eine substituierte Phenanthrenylgruppe, eine substituierte Pyrenylgruppe oder eine substituierte Fluorenylgruppe ist;
R_1 mindestens eine elektronenziehende Gruppe ist, ausgewählt aus einer Pyridinylgruppe, einer Pyrazinylgruppe, einer Pyrimidinylgruppe, einer Chinolinylgruppe, einer Isochinolinylgruppe, einer Chinazolinylgruppe, einer Triazinylgruppe, einer Benzimidazolylgruppe und einer Carbazolylgruppe; und einer Pyridinylgruppe, einer Pyrazinylgruppe, einer Pyrimidinylgruppe, einer Chinolinylgruppe, einer Isochinolinylgruppe, einer Chinazolinylgruppe, einer Phthalazinylgruppe, einer Benzimidazolylgruppe und einer Carbazolylgruppe, die substituiert sind mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN; einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, einer Phosphorsäure oder einem Salz davon, einem C_1-C_{20}-Alkylrest, einem mit mindestens einem -F substituierten C_1-C_{20}-Alkylrest, einem C_1-C_{20}-Alkoxyrest, einer Phenylgruppe, einer Naphthylgruppe, einer Anthrylgruppe, einer Phenanthrenylgruppe, einer Pyrenylgruppe, einer Pyridinylgruppe, einer Triazinylgruppe und einer Carbazolylgruppe; und p 2, 3 oder 4 ist; oder wobei die aminbasierte Verbindung durch die Formel 1(2) wiedergegeben wird, wobei die Ringgruppe A und die Ringgruppe B jeweils unabhängig voneinander eine substituierte Phenylgruppe, eine substituierte Biphenylgruppe, eine substituierte Naphthylgruppe, eine substituierte Anthrylgruppe, eine substituierte Phenanthrenylgruppe, eine substituierte Pyrenylgruppe oder eine substituierte Fluorenylgruppe sind;
R_1 und R_2 jeweils unabhängig voneinander mindestens eine elektronenziehende Gruppe sind, ausgewählt aus einer Pyridinylgruppe, einer Pyrazinylgruppe, einer Pyrimidinylgruppe, einer Chinolinylgruppe, einer Isochinolinylgruppe, einer Chinazolinylgruppe, einer Triazinylgruppe, einer Benzimidazolylgruppe und einer Carbazolylgruppe; und einer

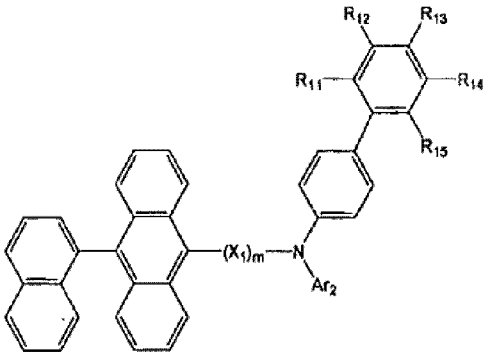
Pyridinylgruppe, einer Pyrazinylgruppe, einer Pyrimidinylgruppe, einer Chinolinylgruppe, einer Isochinolinylgruppe, einer Chinazolinylgruppe, einer Phthalazinylgruppe, einer Benzimidazolylgruppe und einer Carbazolylgruppe, die substituiert sind mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN; einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, einer Phosphorsäure oder einem Salz davon, einem C_1-C_{20}-Alkylrest, einem mit mindestens einem -F substituierten C_1-C_{20}-Alkylrest, einem C_1-C_{20}-Alkoxyrest, einer Phenylgruppe, einer Naphthylgruppe, einer Anthrylgruppe, einer Phenanthrenylgruppe, einer Pyrenylgruppe, einer Pyridinylgruppe, einer Triazinylgruppe und einer Carbazolylgruppe; und p und q jeweils unabhängig voneinander 2, 3 oder 4 sind.

11. Aminbasierte Verbindung nach Anspruch 1, wobei die aminbasierte Verbindung durch eine der nachstehenden Formeln 1A bis 1J wiedergegeben wird:

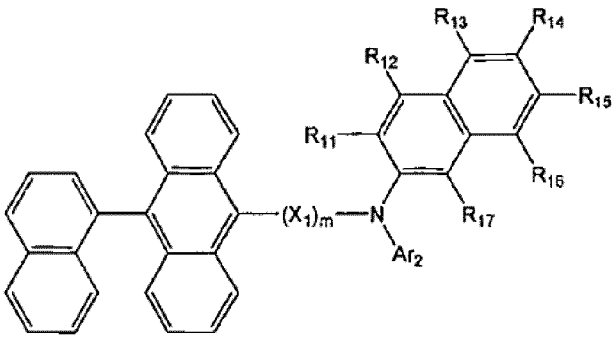
Formel 1A



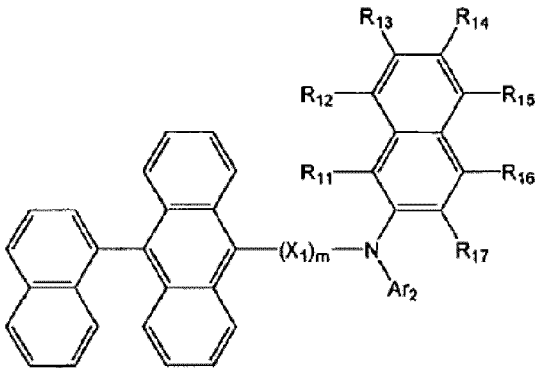
Formel 1B



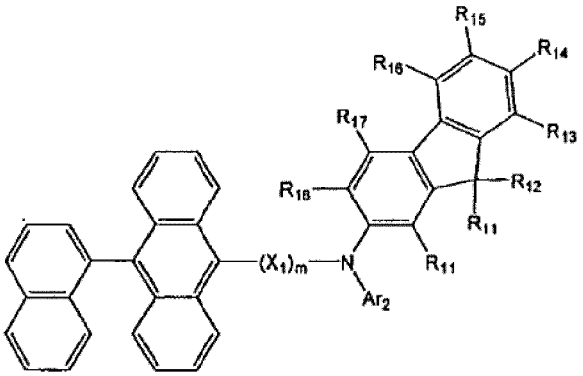
Formel 1C



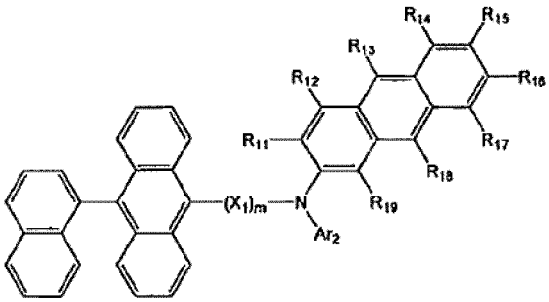
Formel 1D



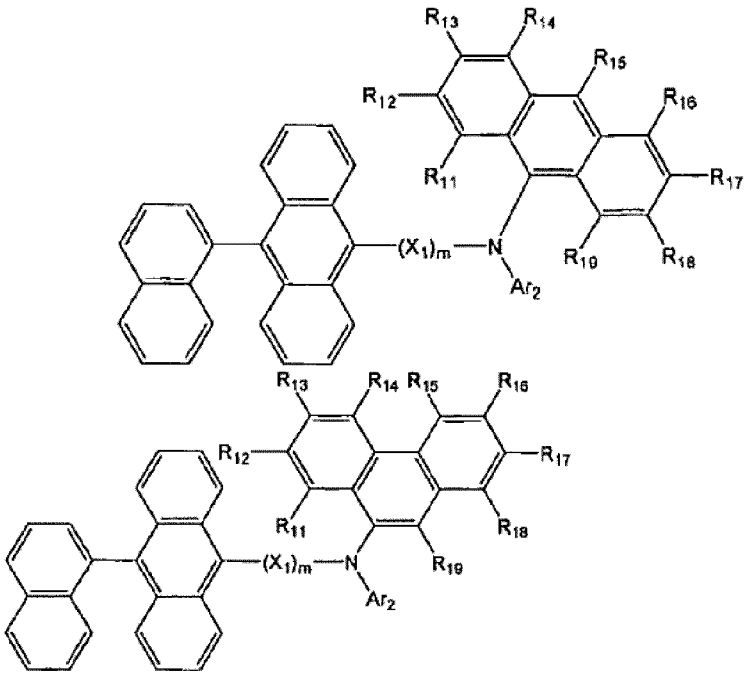
Formel 1E



Formel 1F

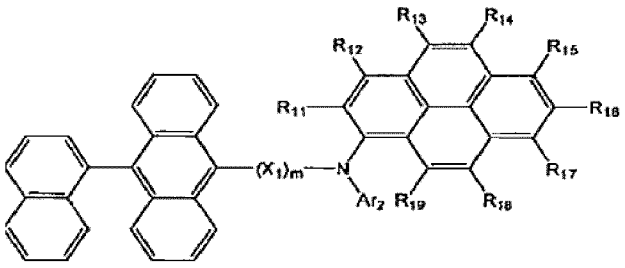


Formel 1G

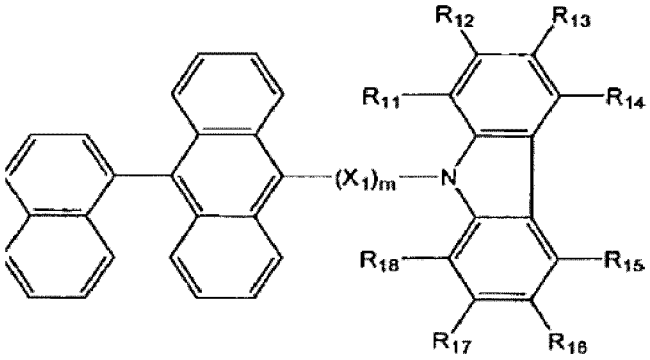


Formel 1H

Formel 1I



Formel 1J



wobei, in den Formeln 1A bis 1J, Ar_2 eine substituierte oder unsubstituierte Phenylgruppe, eine substituierte oder unsubstituierte Pentalenylgruppe, eine substituierte oder unsubstituierte Indenylengruppe, eine substituierte oder unsubstituierte Naphthylgruppe, eine substituierte oder unsubstituierte Azulenylgruppe, eine substituierte oder unsubstituierte Heptalenylgruppe, eine substituierte oder unsubstituierte Indacenylgruppe, eine substituierte oder unsubstituierte Acenaphthylgruppe, eine substituierte oder unsubstituierte Fluorenylgruppe, eine substituierte oder unsubstituierte Phenalenylgruppe, eine substituierte oder unsubstituierte Phenantrhenylgruppe, eine substituierte oder unsubstituierte Anthrylgruppe, eine substituierte oder unsubstituierte Fluoranthenylgruppe, eine substituierte oder unsubstituierte Triphenylenylgruppe, eine substituierte oder unsubstituierte Pyrenylgruppe, eine substituierte oder unsubstituierte Chrysenylgruppe, eine substituierte oder unsubstituierte Naphthacenylgruppe, eine substituierte oder unsubstituierte Picenylgruppe, eine substituierte oder unsubstituierte Perylenylgruppe, eine substituierte oder unsubstituierte Pentaphenylgruppe, eine substituierte oder unsubstituierte Hexacenylgruppe, eine substituierte oder unsubstituierte Pyrrolylgruppe, eine substituierte oder unsubstituierte Pyrazolylgruppe, eine substituierte oder unsubstituierte Imidazolylgruppe, eine substituierte oder unsubstituierte Imidazolinylgruppe, eine substituierte oder unsubstituierte Imidazopyridinylgruppe, eine substituierte oder unsubstituierte Imidazopyrimidinylgruppe, eine substituierte oder unsubstituierte Pyridinylgruppe, eine substituierte oder unsubstituierte Pyrazinylgruppe, eine substituierte oder unsubstituierte Pyrimidinylgruppe, eine substituierte oder unsubstituierte Benzimidazolylgruppe, eine substituierte oder unsubstituierte Indolylgruppe, eine substituierte oder unsubstituierte Purinylgruppe, eine substituierte oder unsubstituierte Chinolinylgruppe, eine substituierte oder unsubstituierte Phthalazinylgruppe, eine substituierte oder unsubstituierte Indolizinylgruppe, eine substituierte oder unsubstituierte Naphthyridinylgruppe, eine substituierte oder unsubstituierte Chinazolinylgruppe, eine substituierte oder unsubstituierte Cinnolinylgruppe, eine substituierte oder unsubstituierte Indazolylgruppe, eine substituierte oder unsubstituierte Carbazolylgruppe, eine substituierte oder unsubstituierte Phenazinylgruppe, eine substituierte oder unsubstituierte Phenanthridinylgruppe, eine substituierte oder unsubstituierte Pyranylgruppe, eine substituierte oder unsubstituierte Chromenylgruppe, eine substituierte oder unsubstituierte Furanylgruppe, eine substituierte oder unsubstituierte Benzofuranylgruppe, eine substituierte oder unsubstituierte Thiophenylgruppe, eine substituierte oder unsubstituierte Benzothiophenylgruppe, eine substituierte oder unsubstituierte Isothiazolylgruppe, eine substituierte oder unsubstituierte Benzimidazolylgruppe, eine substituierte oder unsubstituierte Isoxazolylgruppe, eine substituierte oder unsubstituierte Dibenzothiophenylgruppe, eine substituierte oder unsubstituierte Dibenzofuranylgruppe, eine substituierte oder unsubstituierte Triazinylgruppe, eine substituierte oder unsubstituierte Oxadiazolylgruppe, eine substituierte oder unsubstituierte Pyridazinylgruppe, eine substituierte oder unsubstituierte Triazolylgruppe, eine substituierte oder unsubstituierte Tetrazolylgruppe oder eine substituierte oder unsubstituierte Phenanthrolinylgruppe ist,

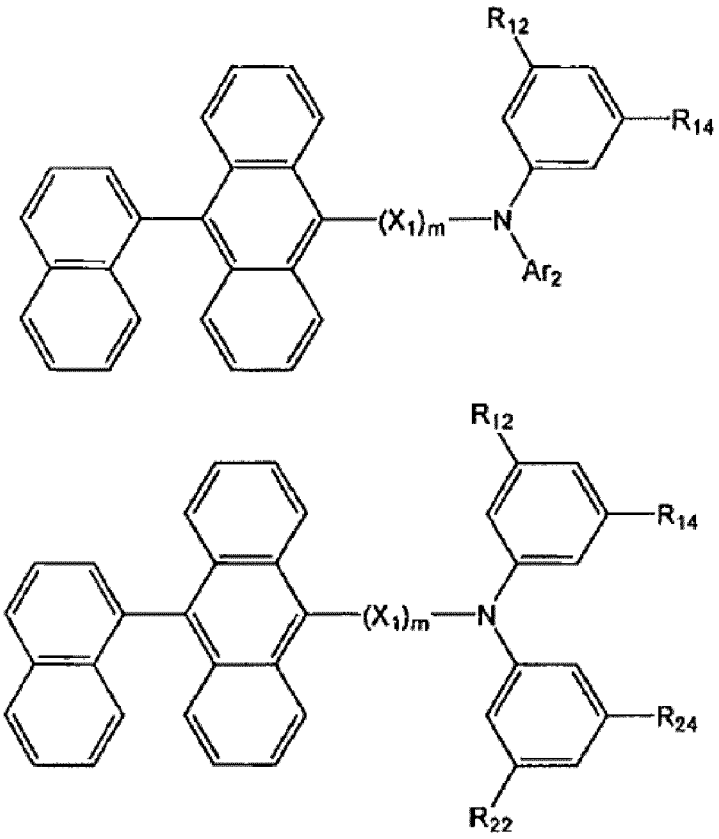
Substituenten der substituierten Phenylgruppe, der substituierten Pentalenylgruppe, der substituierten Indenylgruppe, der substituierten Naphthylgruppe, der substituierten Azulenylgruppe, der substituierten Heptalenylgruppe, der

substituierten Indacenylgruppe, der substituierten Acenaphthylgruppe, der substituierten Fluorenylgruppe, der substituierten Phenalenylgruppe, der substituierten Phenanthrenylgruppe, der substituierten Anthrylgruppe, der substituierten Fluoranthenylgruppe, der substituierten Triphenylenylgruppe, der substituierten Pyrenylgruppe, der substituierten Chrysenylgruppe, der substituierten Naphthacenylgruppe, der substituierten Picenylgruppe, der substituierten Perylenylgruppe, der substituierten Pentaphenylgruppe, der substituierten Hexacenylgruppe, der substituierten Pyrrolylgruppe, der substituierten Pyrazolylgruppe, der substituierten Imidazolylgruppe, der substituierten Imidazolinylgruppe, der substituierten Imidazopyridinylgruppe, der substituierten Imidazopyrimidinylgruppe, der substituierten Pyridinylgruppe, der substituierten Pyrazinylgruppe, der substituierten Pyrimidinylgruppe, der substituierten Benzimidazolylgruppe, der substituierten Indolylgruppe, der substituierten Purinylgruppe, der substituierten Chinolinylgruppe, der substituierten Phthalazinylgruppe, der substituierten Indolizinylgruppe, der substituierten Naphthyridinylgruppe, der substituierten Chinazolinylgruppe, der substituierten Cinnolinylgruppe, der substituierten Indazolylgruppe, der substituierten Carbazolylgruppe, der substituierten Phenazinylgruppe, der substituierten Phenanthridinylgruppe, der substituierten Pyranylgruppe, der substituierten Chromenylgruppe, der substituierten Furanylgruppe, der substituierten Benzofuranylgruppe, der substituierten Thiophenylgruppe, der substituierten Benzothiophenylgruppe, der substituierten Isothiazolylgruppe, der substituierten Benzimidazolylgruppe, der substituierten Isoxazolylgruppe, der substituierten Dibenzothiophenylgruppe, der substituierten Dibenzofuranylgruppe, der substituierten Triazinylgruppe, der substituierten Oxadiazolylgruppe, der substituierten Pyridazinylgruppe, der substituierten Triazolylgruppe, der substituierten Tetrazolylgruppe und der substituierten Phenanthrolinylgruppe, und R_{11} bis R_{19} jeweils unabhängig voneinander ein Wasserstoffatom; ein Deuteriumatom; -F; -Cl; -Br; -I; -CN; eine Hydroxylgruppe; $-NO_2$; eine Aminogruppe; eine Amidinogruppe; Hydrazin; Hydrazon; eine Carboxylgruppe oder ein Salz davon; eine Sulfonsäuregruppe oder ein Salz davon; Phosphorsäure oder ein Salz davon; ein Tri(C_6-C_{60}-aryl)silylrest; ein C_1-C_{60}-Alkylrest, ein C_1-C_{60}-Alkoxyrest, ein C_2-C_{60}-Alkenylrest und ein C_2-C_{60}-Alkinylrest; ein C_1-C_{60}-Alkylrest, ein C_1-C_{60}-Alkoxyrest, ein C_2-C_{60}-Alkenylrest und ein C_2-C_{60}-Alkinylrest, die substituiert sind mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon und Phosphorsäure oder einem Salz davon; ein C_3-C_{60}-Cycloalkylrest, ein C_3-C_{60}-Cycloalkenylrest, ein C_6-C_{60}-Arylrest, ein C_2-C_{60}-Heteroarylrest, ein C_6-C_{60}-Aralkylrest, ein C_6-C_{60}-Aryloxyrest und ein C_6-C_{60}-Arylthiolrest; und ein C_3-C_{60}-Cycloalkylrest, ein C_3-C_{60}-Cycloalkenylrest, ein C_6-C_{60}-Arylrest, ein C_2-C_{60}-Heteroarylrest, ein C_6-C_{60}-Aralkylrest, ein C_6-C_{60}-Aryloxyrest und ein C_6-C_{60}-Arylthiolrest, die substituiert sind mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, Phosphorsäure oder einem Salz davon, einem C_1-C_{60}-Alkylrest, einem mit mindestens einem -F substituierten C_1-C_{60}-Alkylrest, einem C_1-C_{60}-Alkoxyrest, einem C_2-C_{60}-Alkenylrest, einem C_2-C_{60}-Alkinylrest; einem C_6-C_{60}-Arylrest und einem C_2-C_{60}-Heteroarylrest, sind,

wobei mindestens eines von R_{11} bis R_{15} in den Formeln 1A und 1B, mindestens eines von R_{11} bis R_{17} in den Formeln 1C und 1D, mindestens eines von R_{11} bis R_{18} in den Formeln 1E und 1J und mindestens eines von R_{11} bis R_{19} in Formel 1F, 1G, 1H und 1I jeweils unabhängig voneinander eine elektronenziehende Gruppe sind, ausgewählt aus -F; -CN; $-NO_2$, einem mit mindestens einem -F substituierten C_1-C_{60}-Alkylrest; einem C_2-C_{60}-Heteroarylrest; und einem C_2-C_{60}-Heteroarylrest, substituiert mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, einer Phosphorsäure oder einem Salz davon, einem C_1-C_{60}-Alkylrest, einem mit mindestens einem -F substituierten C_1-C_{60}-Alkylrest, einem C_1-C_{60}-Alkoxyrest, einem C_2-C_{60}-Alkenylrest, einem C_2-C_{60}-Alkinylrest; einem C_6-C_{60}-Arylrest und einem C_2-C_{60}-Heteroarylrest.

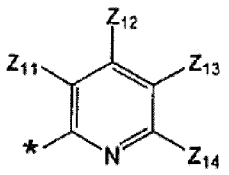
12. Aminbasierte Verbindung nach Anspruch 11, wobei die aminbasierte Verbindung durch die nachstehende Formel 1A-(1) oder 1A-(2) wiedergegeben wird:

Formel 1A-(1) Formel 1A-(2)

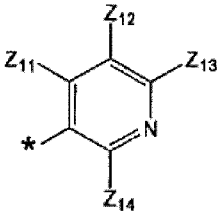


wobei, in den Formeln 1A-(1) oder 1A-(2), R_{12}, R_{14}, R_{22} und R_{24} jeweils unabhängig voneinander eine elektronenziehende Gruppe sind, ausgewählt aus einer Pyridinylgruppe, einer Pyrazinylgruppe, einer Pyrimidinylgruppe, einer Chinolinylgruppe, einer Isochinolinylgruppe, einer Chinazolinylgruppe, einer Triazinylgruppe, einer Benzimidazolylgruppe und einer Carbazolylgruppe; und einer Pyridinylgruppe, einer Pyrazinylgruppe, einer Pyrimidinylgruppe, einer Chinolinylgruppe, einer Isochinolinylgruppe, einer Chinazolinylgruppe, einer Phthalazinylgruppe, einer Benzimidazolylgruppe und einer Carbazolylgruppe, die substituiert sind mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN; einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, einer Phosphorsäure oder einem Salz davon, einem C_1-C_{20}-Alkylrest, einem mit mindestens einem -F substituierten C_1-C_{20}-Alkylrest, einem C_1-C_{20}-Alkoxyrest, einer Phenylgruppe, einer Naphthylgruppe, einer Anthrylgruppe, einer Phenanthrenylgruppe, einer Pyrenylgruppe, einer Pyridinylgruppe, einer Triazinylgruppe und einer Carbazolylgruppe; und
in Formel 1A-(1) Ar_2 eine substituierte oder unsubstituierte Phenylgruppe, eine substituierte oder unsubstituierte Biphenylgruppe, eine substituierte oder unsubstituierte Naphthylgruppe, eine substituierte oder unsubstituierte Anthrylgruppe, eine substituierte oder unsubstituierte Phenanthrenylgruppe, eine substituierte oder unsubstituierte Pyrenylgruppe oder eine substituierte oder unsubstituierte Fluorenylgruppe ist.

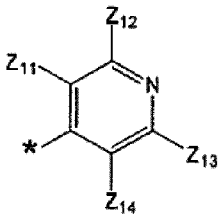
13. Aminbasierte Verbindung nach Anspruch 12, wobei R_{12}, R_{14}, R_{22} und R_{24} jeweils unabhängig voneinander ausgewählt sind aus F; -CN; $-CH_2F$, $-CHF_2$; $-CF_3$ und den durch die nachstehenden Formeln 2(1) bis 2(14) dargestellten Resten:



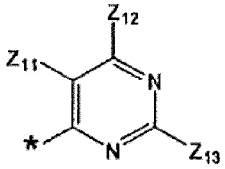
Formel 2(1)



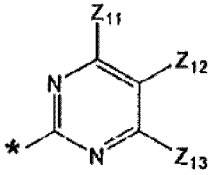
Formel 2(2)



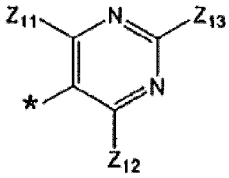
Formel 2(3)



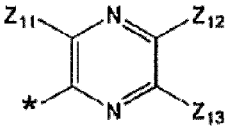
Formel 2(4)



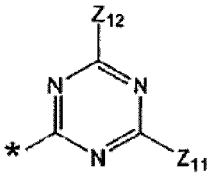
Formel 2(5)



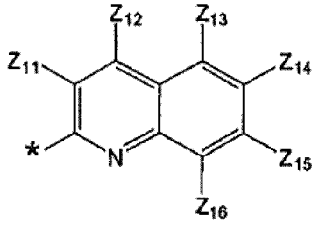
Formel 2(6)



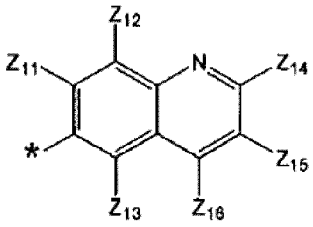
Formel 2(7)



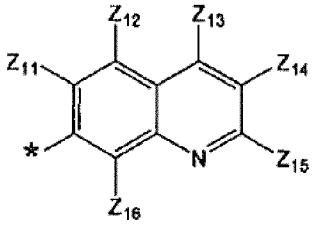
Formel 2(8)



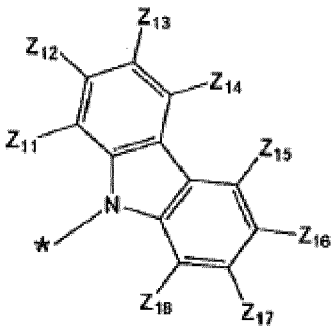
Formel 2(9)



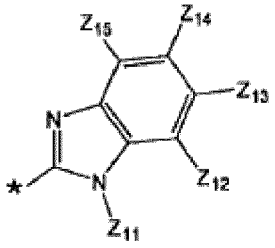
Formel 2(10)



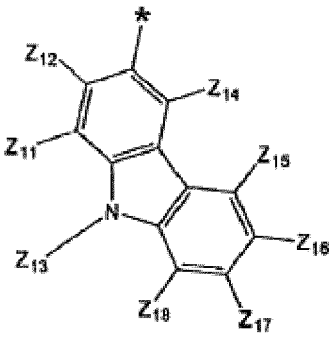
Formel 2(11)



Formel 2(12)



Formel 2(13)



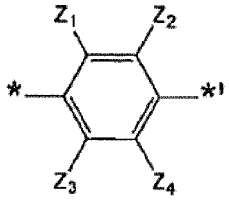
Formel 2(14)

wobei, in den Formeln 2(1) bis 2(14), Z_{11} bis Z_{18} jeweils unabhängig voneinander ein Wasserstoffatom, ein Deuteriumatom, -F, -Cl, -Br, -I, -CN, eine Hydroxylgruppe, $-NO_2$, eine Aminogruppe, eine Amidinogruppe, Hydrazin, Hydrazon, eine Carboxylgruppe oder ein Salz davon, eine Sulfonsäuregruppe oder ein Salz davon, Phosphorsäure oder ein Salz davon, ein C_1-C_{20}-Alkylrest, ein mit mindestens einem -F substituierter C_1-C_{20}-Alkylrest, ein C_1-C_{20}-Alkoxyrest, eine Phenylgruppe, eine Naphthylgruppe, eine Anthrylgruppe, eine Phenanthrenylgruppe, eine Pyrenylgruppe, eine Pyridinylgruppe, eine Triazinylgruppe oder eine Carbazolylgruppe sind.

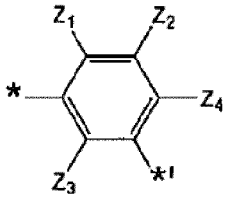
14. Aminbasierte Verbindung nach einem vorangehenden Anspruch, wobei X_1 eine substituierte oder unsubstituierte Phenylengruppe, eine substituierte oder unsubstituierte Pentalenylengruppe, eine substituierte oder unsubstituierte Indenylengruppe, eine substituierte oder unsubstituierte Naphthylengruppe, eine substituierte oder unsubstituierte Azulenylengruppe, eine substituierte oder unsubstituierte Heptalenylengruppe, eine substituierte oder unsubstituierte Indacenylengruppe, eine substituierte oder unsubstituierte Acenaphthylengruppe, eine substituierte oder un-

substituierte Fluorenylengruppe, eine substituierte oder unsubstituierte Phenalenylengruppe, eine substituierte oder unsubstituierte Phenanthrenylengruppe, eine substituierte oder unsubstituierte Anthrylengruppe, eine substituierte oder unsubstituierte Fluoranthenylengruppe, eine substituierte oder unsubstituierte Triphenylenylengruppe, eine substituierte oder unsubstituierte Pyrenylengruppe, eine substituierte oder unsubstituierte Chrysenylengruppe, eine substituierte oder unsubstituierte Naphthacenylengruppe eine substituierte oder unsubstituierte Picenylengruppe, eine substituierte oder unsubstituierte Perylenylengruppe, eine substituierte oder unsubstituierte Pentaphenylengruppe, eine substituierte oder unsubstituierte Hexacenylengruppe, eine substituierte oder unsubstituierte Pyrrolylengruppe, eine substituierte oder unsubstituierte Pyrazolylengruppe, eine substituierte oder unsubstituierte Imidazolylengruppe, eine substituierte oder unsubstituierte Imidazolinylengruppe, eine substituierte oder unsubstituierte Imidazopyridinylengruppe, eine substituierte oder unsubstituierte Imidazopyrimidinylengruppe, eine substituierte oder unsubstituierte Pyridinylengruppe, eine substituierte oder unsubstituierte Pyrazinylengruppe, eine substituierte oder unsubstituierte Pyrimidinylengruppe, eine substituierte oder unsubstituierte Indolylengruppe, eine substituierte oder unsubstituierte Purinylengruppe, eine substituierte oder unsubstituierte Chinolinylengruppe, eine substituierte oder unsubstituierte Phthalazinylengruppe, eine substituierte oder unsubstituierte Indolizinylengruppe, eine substituierte oder unsubstituierte Naphthyridinylengruppe, eine substituierte oder unsubstituierte Chinazolinylengruppe, eine substituierte oder unsubstituierte Cinnolinylengruppe, eine substituierte oder unsubstituierte Indazolylengruppe, eine substituierte oder unsubstituierte Carbazolylengruppe, eine substituierte oder unsubstituierte Phenazinylengruppe, eine substituierte oder unsubstituierte Phenanthridinylengruppe, eine substituierte oder unsubstituierte Pyranylengruppe, eine substituierte oder unsubstituierte Chromenylengruppe, eine substituierte oder unsubstituierte Furanylengruppe, eine substituierte oder unsubstituierte Benzofuranylengruppe, eine substituierte oder unsubstituierte Thiophenylengruppe, eine substituierte oder unsubstituierte Benzothiophenylengruppe, eine substituierte oder unsubstituierte Isothiazolylengruppe, eine substituierte oder unsubstituierte Benzimidazolylengruppe, eine substituierte oder unsubstituierte Isoxazolylengruppe, eine substituierte oder unsubstituierte Dibenzothiophenylengruppe, eine substituierte oder unsubstituierte Dibenzofuranylengruppe, eine substituierte oder unsubstituierte Triazinylengruppe, eine substituierte oder unsubstituierte Oxadiazolylengruppe, eine substituierte oder unsubstituierte Pyridazinylengruppe, eine substituierte oder unsubstituierte Triazolylengruppe oder eine substituierte oder unsubstituierte Tetrazolylengruppe ist.

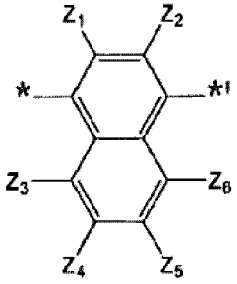
15. Aminbasierte Verbindung nach Anspruch 14, wobei X_1 ein durch eine der nachstehenden Formeln 5(1) bis 5(16) dargestellter Rest ist:



Formel 5(1)



Formel 5(2)



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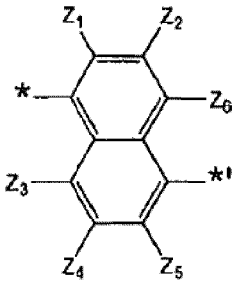

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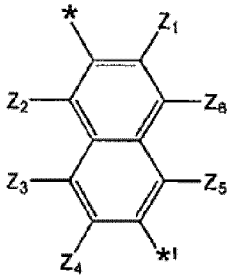

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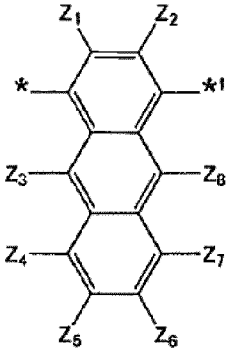

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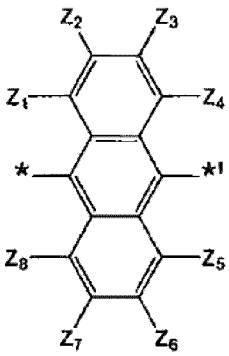

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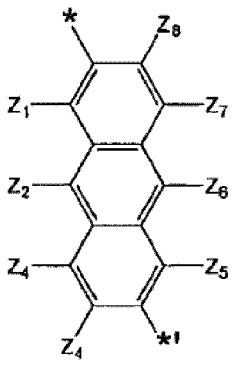

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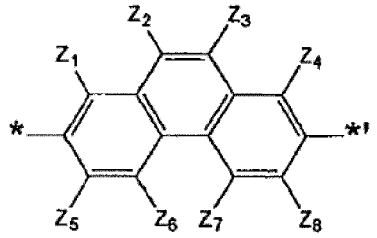

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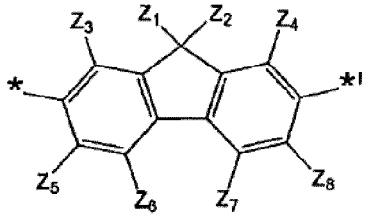

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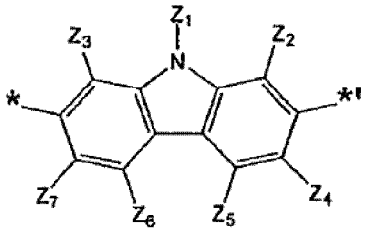

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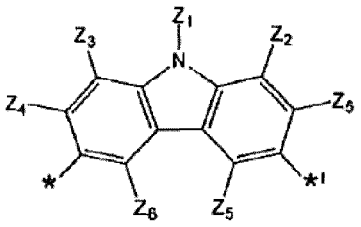

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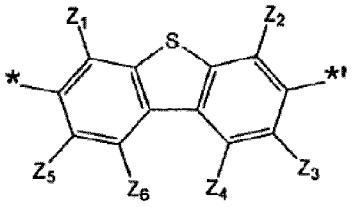
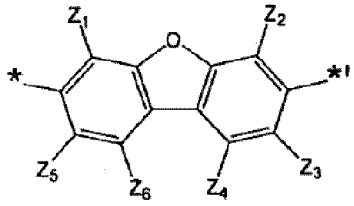


Formel 5(13)



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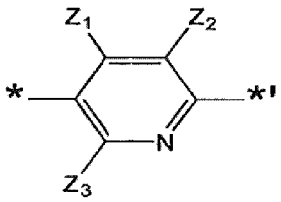
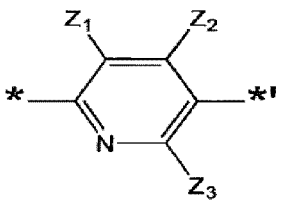


Formel 5(15)



Formel 5(16)

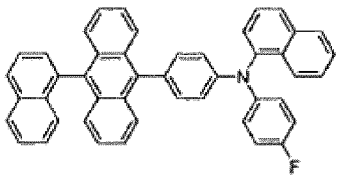
wobei, in den Formeln 5(1) bis 5(16), Z_1 bis Z_8 jeweils unabhängig voneinander eines von einem Wasserstoffatom; einem Deuteriumatom; -F; -Cl; -Br; -I; -CN; einer Hydroxylgruppe; $-NO_2$; einer Aminogruppe; einer Amidinogruppe; Hydrazin; Hydrazon; einer Carboxylgruppe oder einem Salz davon; einer Sulfonsäuregruppe oder einem Salz davon; Phosphorsäure oder einem Salz davon; einem C_1-C_{20}-Alkylrest; einem C_1-C_{20}-Alkoxyrest; einem C_1-C_{20}-Alkylrest und einem C_1-C_{20}-Alkoxyrest, die substituiert sind mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon und Phosphorsäure oder einem Salz davon; einem C_6-C_{20}-Arylrest; einem C_2-C_{20}-Heteroarylrest; und einem C_6-C_{20}-Arylrest und einem C_2-C_{20}-Heteroarylrest, die substituiert sind mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, Phosphorsäure oder einem Salz davon, einem C_1-C_{20}-Alkylrest, einem C_1-C_{20}-Alkoxyrest, einem C_6-C_{20}-Arylrest und einem C_2-C_{20}-Heteroarylrest, sind;
wobei * eine Bindungsstelle an Anthracen in Formel 1 kennzeichnet; und
wobei *' eine Bindungsstelle an N in Formel 1 kennzeichnet.

16. Aminbasierte Verbindung nach Anspruch 15, wobei Z_1 bis Z_8 jeweils unabhängig voneinander eines von einem Wasserstoffatom; einem Deuteriumatom; -F; -Cl; -Br; -I; -CN; einer Hydroxylgruppe; $-NO_2$; einer Aminogruppe; einer Amidinogruppe; Hydrazin; Hydrazon; einer Carboxylgruppe oder einem Salz davon; einer Sulfonsäuregruppe oder einem Salz davon; einer Phosphorsäure oder einem Salz davon; einer Methylgruppe, einer Ethylgruppe, einer Propylgruppe, einer Butylgruppe und einer Pentylgruppe; einer Methoxygruppe, einer Ethoxygruppe, einer Propoxygruppe, einer Butoxygruppe und einer Pentoxygruppe; einer Methylgruppe, einer Ethylgruppe, einer Propylgruppe, einer Butylgruppe, einer Pentylgruppe, einer Methoxygruppe, einer Ethoxygruppe, einer Propoxygruppe, einer Butoxygruppe und einer Pentoxygruppe, die substituiert sind mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon und einer Phosphorsäure oder einem Salz davon; einer Phenylgruppe, einer Naphthylgruppe, einer Anthrylgruppe, einer Phenanthrenylgrup-

pe, einer Pyrenylgruppe und einer Fluorenylgruppe; einer Pyridinylgruppe, einer Pyrimidinylgruppe, einer Triazinylgruppe, einer Chinolylgruppe und einer Carbazolylgruppe; einer Phenylgruppe, einer Naphthylgruppe, einer Anthrylgruppe, einer Phenanthrenylgruppe, einer Pyrenylgruppe, einer Fluorenylgruppe; einer Pyridinylgruppe, einer Pyrimidinylgruppe, einer Triazinylgruppe, einer Chinolylgruppe und einer Carbazolylgruppe, die substituiert sind mit mindestens einem von einem Deuteriumatom, -F, -Cl, -Br, -I, -CN, einer Hydroxylgruppe, $-NO_2$, einer Aminogruppe, einer Amidinogruppe, Hydrazin, Hydrazon, einer Carboxylgruppe oder einem Salz davon, einer Sulfonsäuregruppe oder einem Salz davon, einer Phosphorsäure oder einem Salz davon, einem C_1-C_{20}-Alkylrest und einem C_1-C_{20}-Alkoxyrest, sind.

17. Aminbasierte Verbindung nach einem vorangehenden Anspruch, wobei m 1, 2 oder 3 ist.

18. Aminbasierte Verbindung nach einem vorangehenden Anspruch, wobei die aminbasierte Verbindung eine der nachstehenden Verbindungen 1 bis 109 ist:



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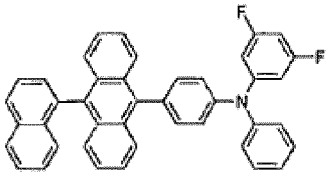

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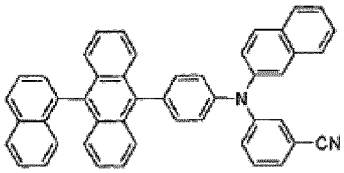

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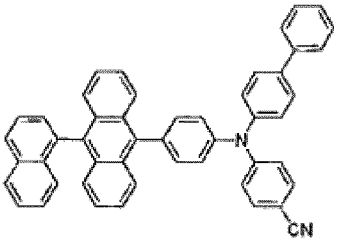

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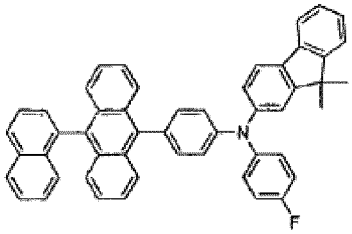

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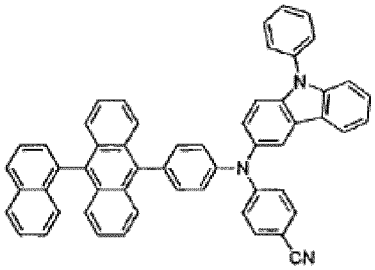

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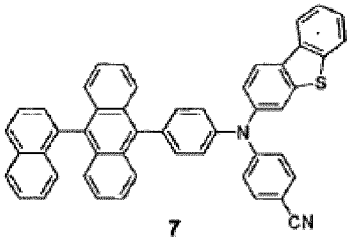

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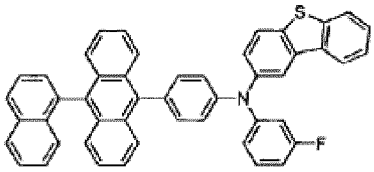

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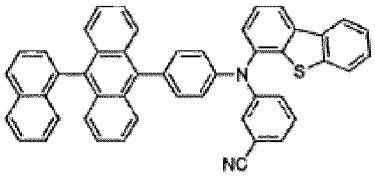

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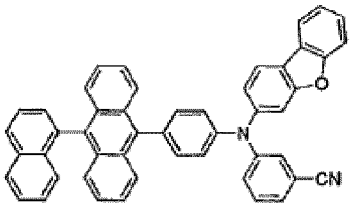

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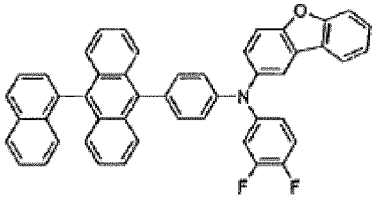

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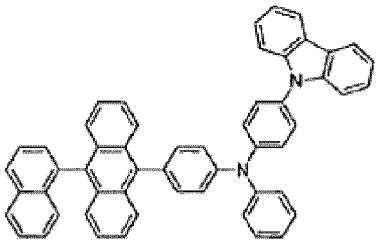

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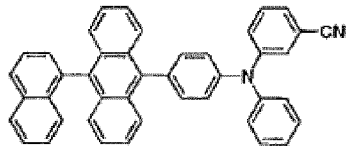

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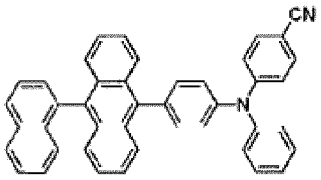

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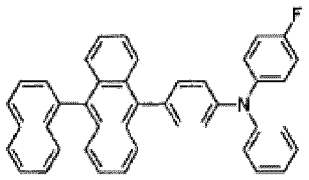

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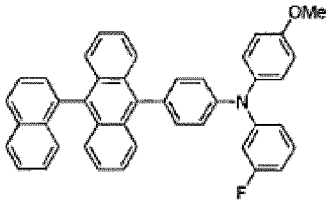

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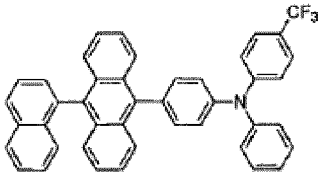

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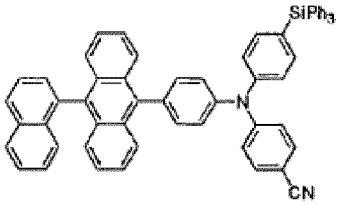

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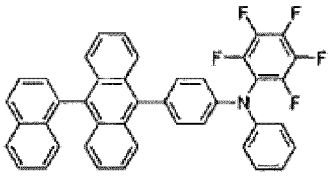

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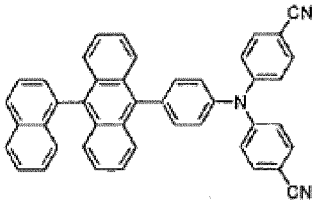

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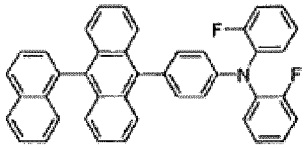

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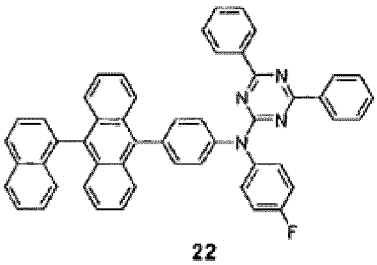

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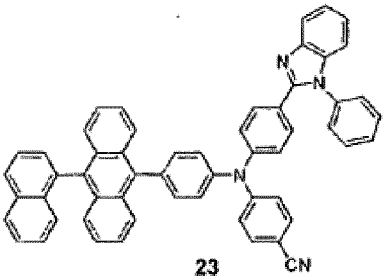

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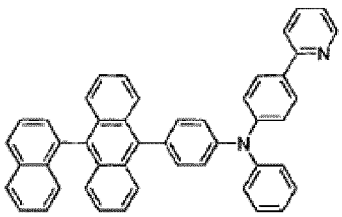

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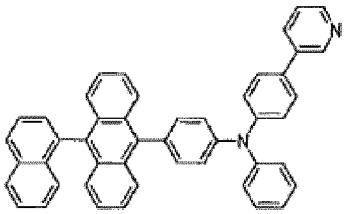

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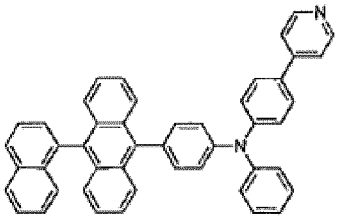

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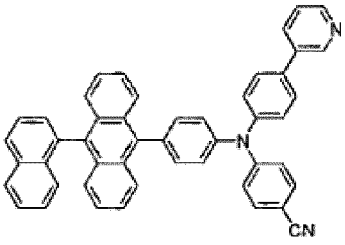

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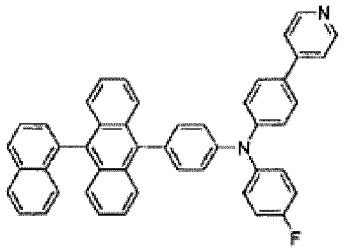

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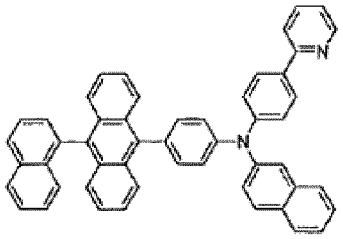

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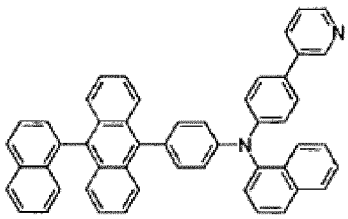

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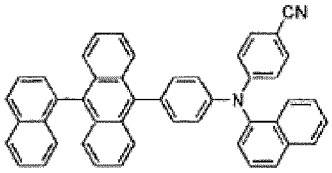

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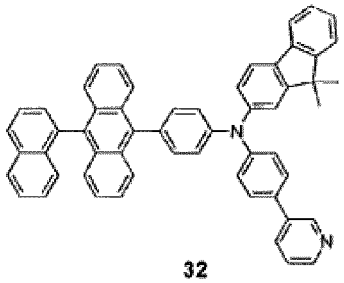

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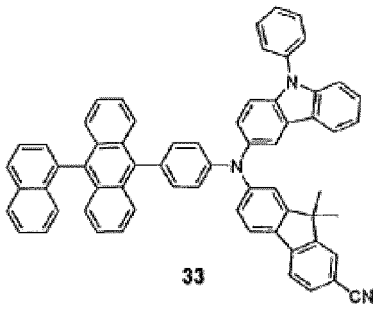

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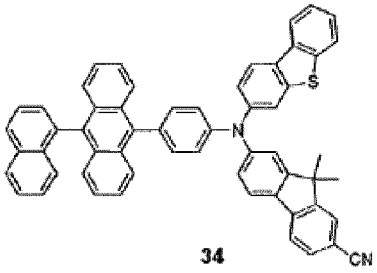

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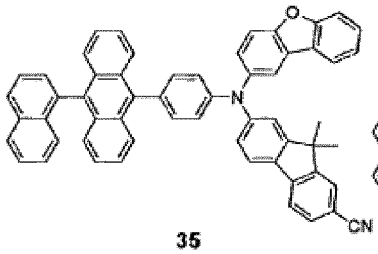

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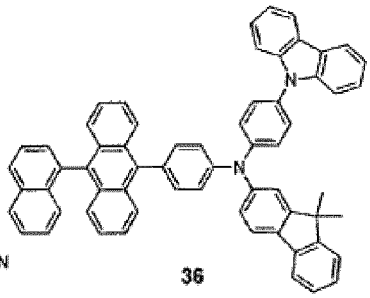

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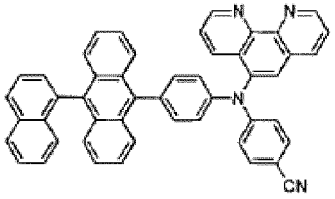

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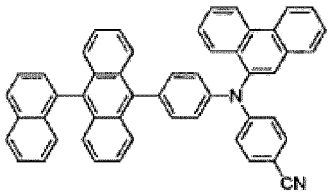

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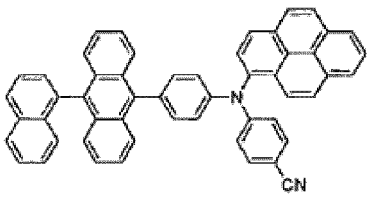

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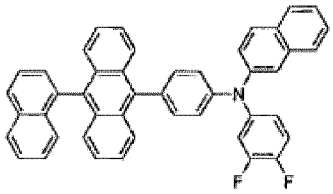

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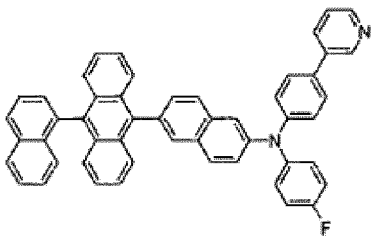

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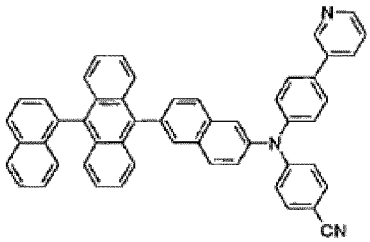

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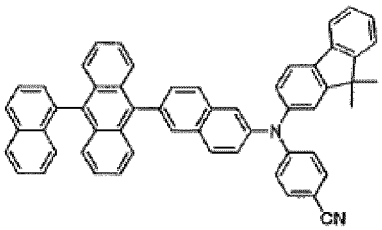

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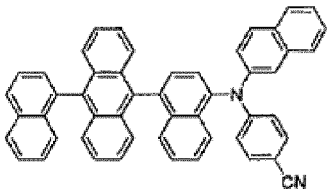

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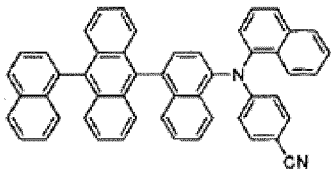

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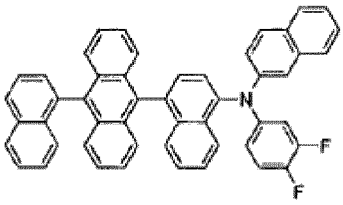

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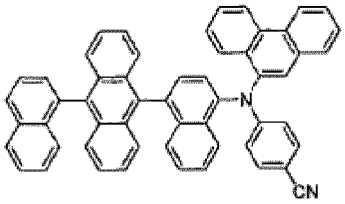

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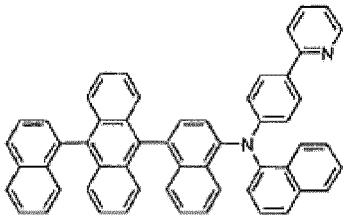

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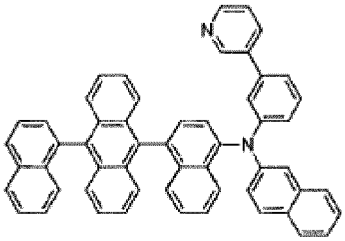

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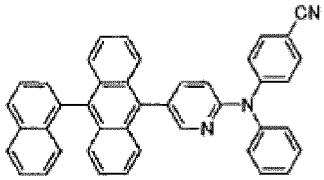

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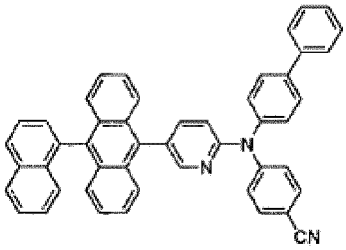

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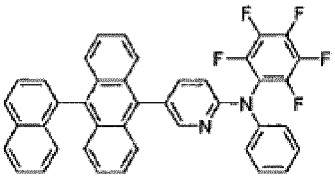

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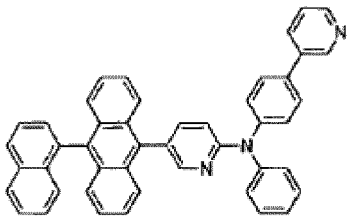

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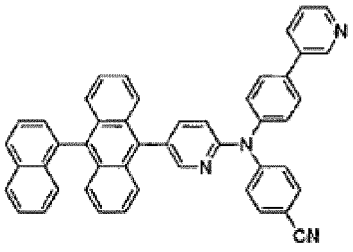

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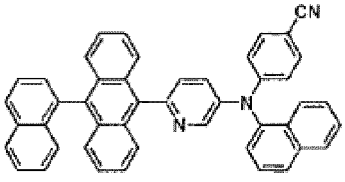

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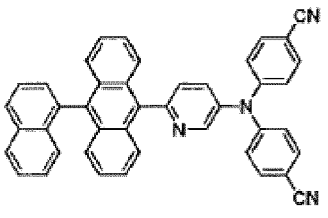

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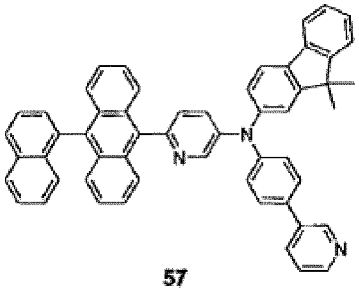

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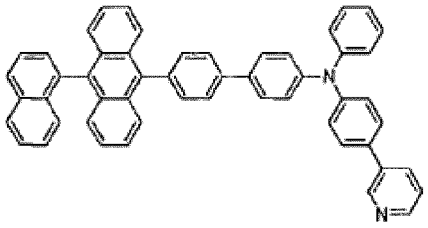

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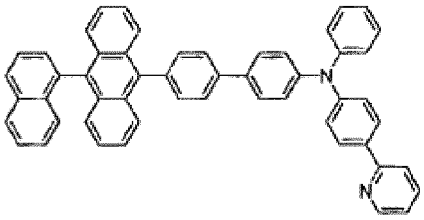

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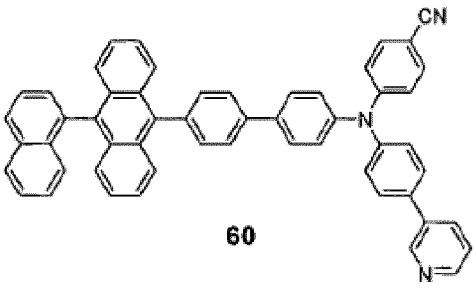

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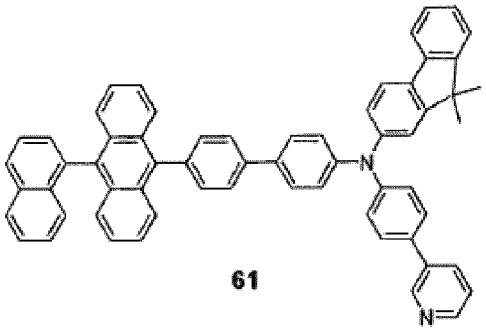

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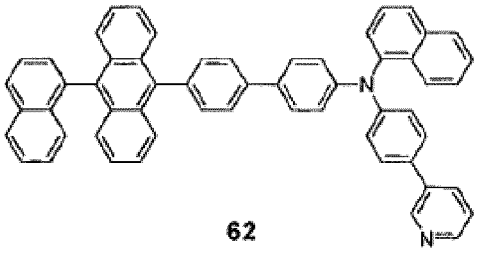

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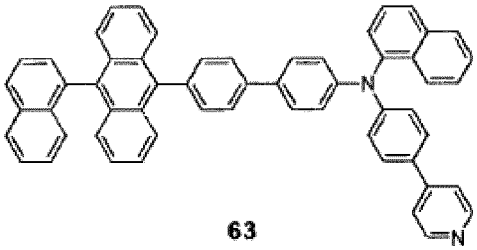

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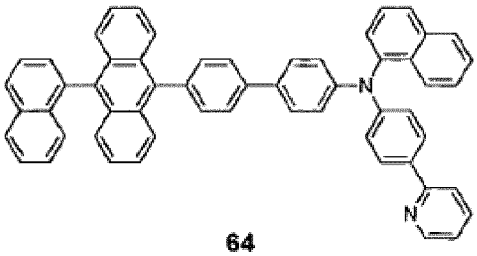

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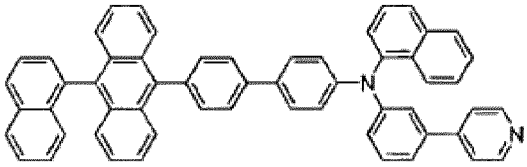

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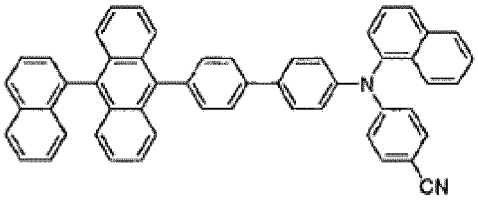

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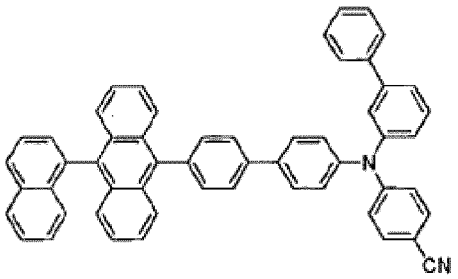

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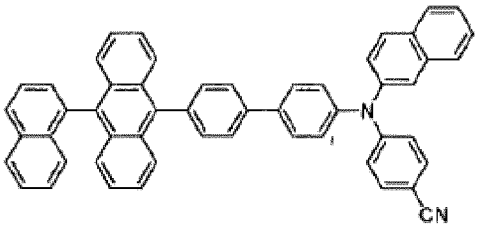

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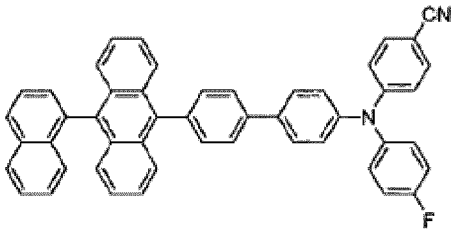

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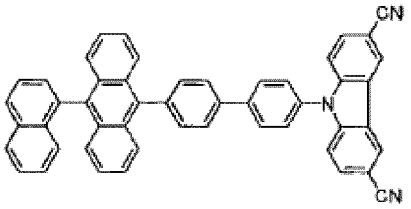

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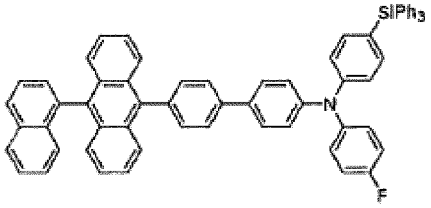

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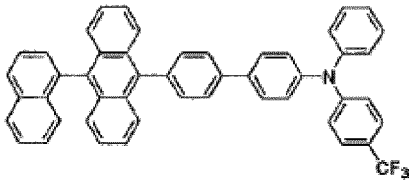

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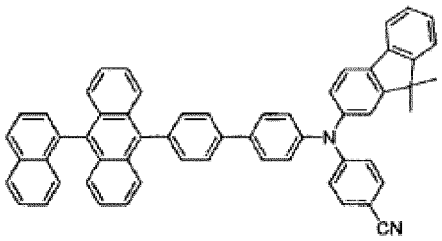

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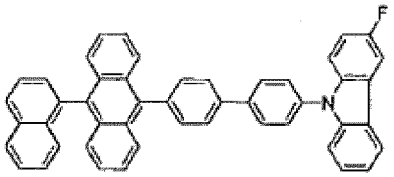

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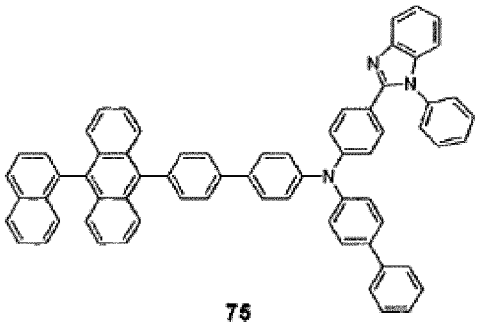

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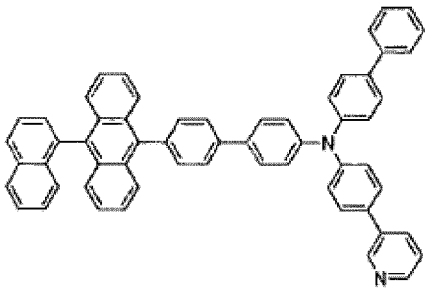

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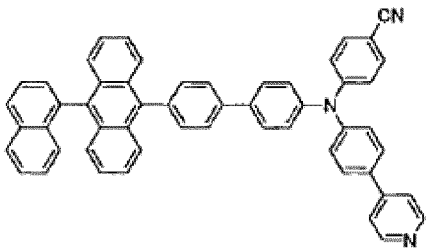

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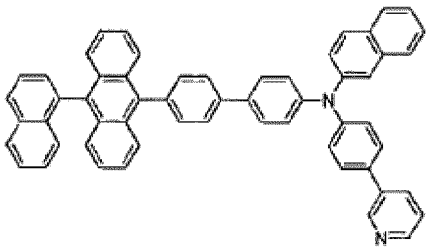

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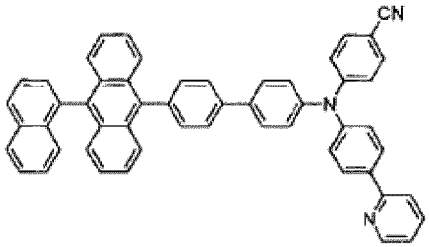

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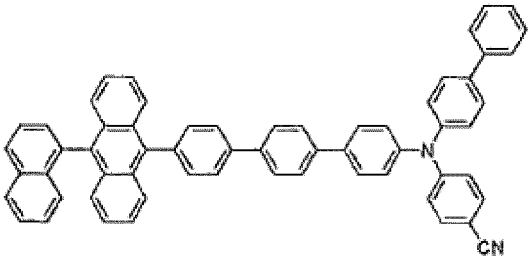

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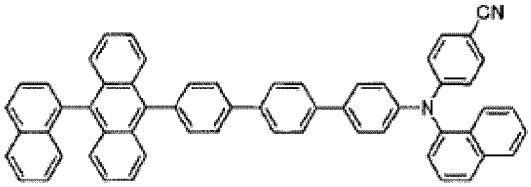

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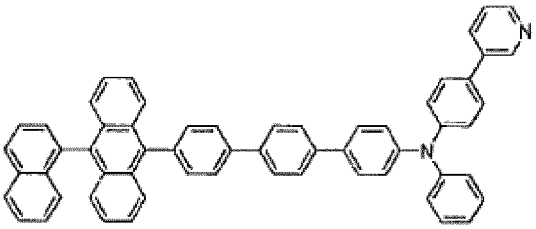

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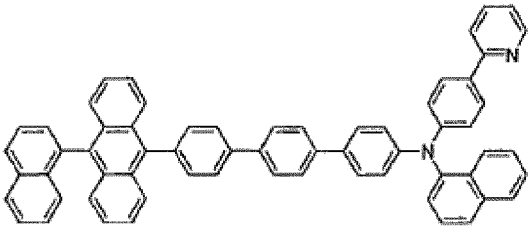

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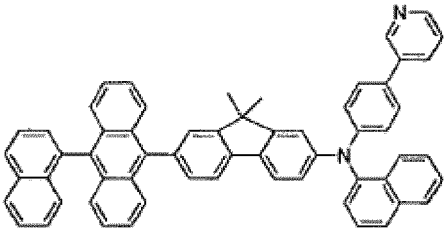

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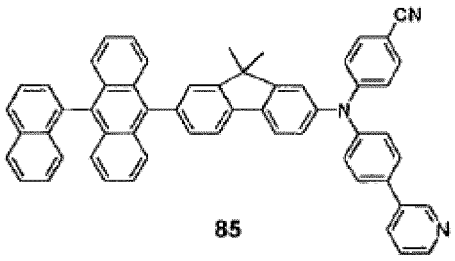

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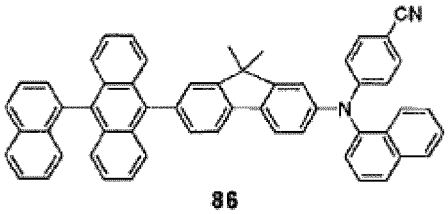

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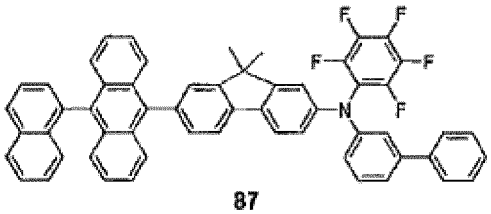

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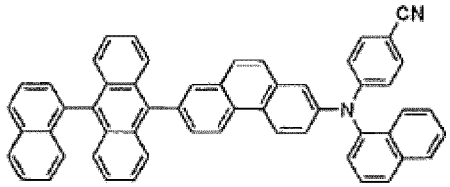

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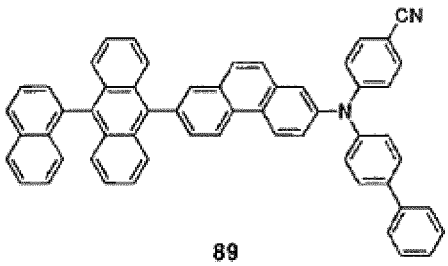

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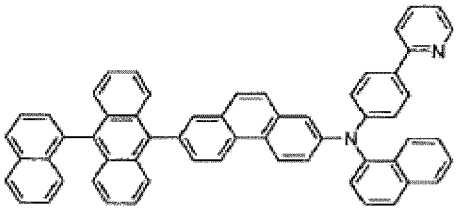

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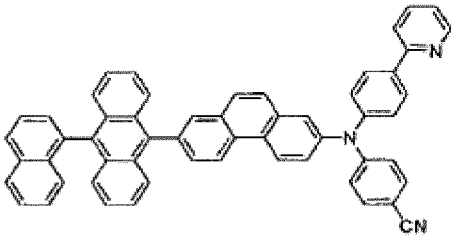

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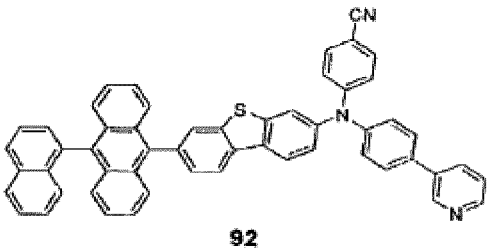

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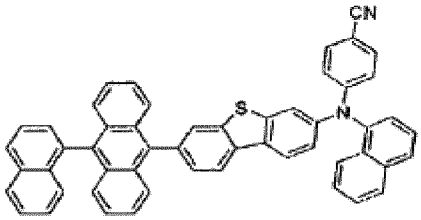

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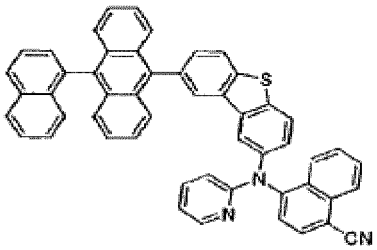

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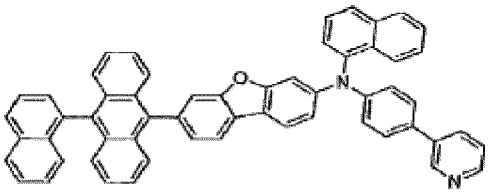

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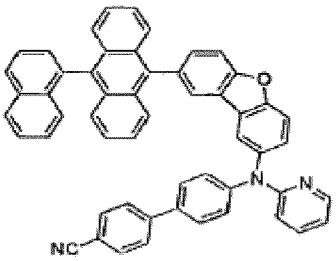

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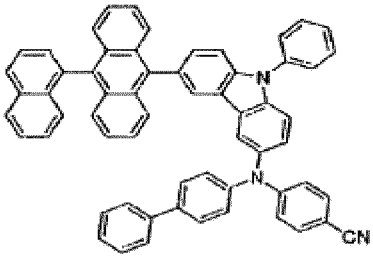

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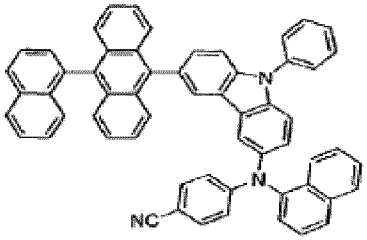

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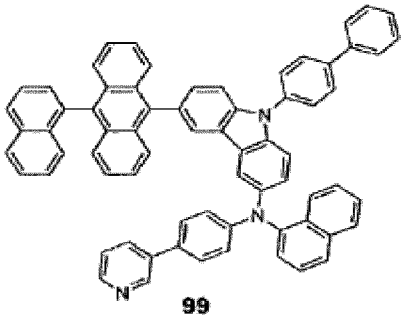

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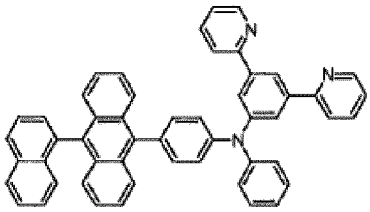

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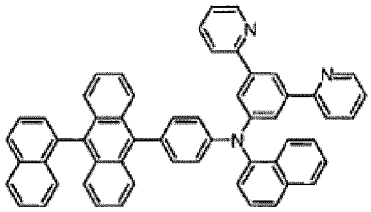

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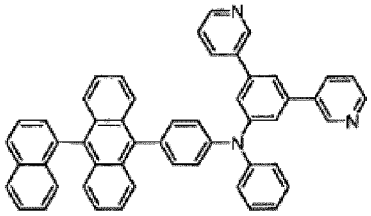

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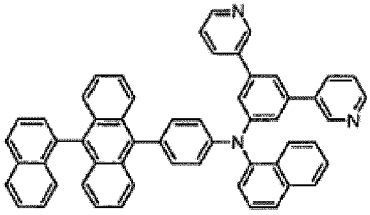

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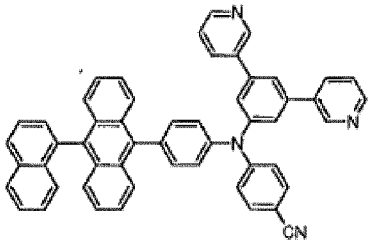

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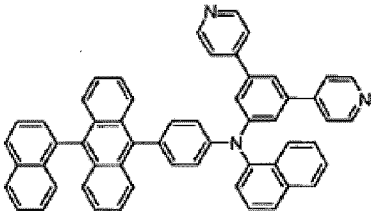

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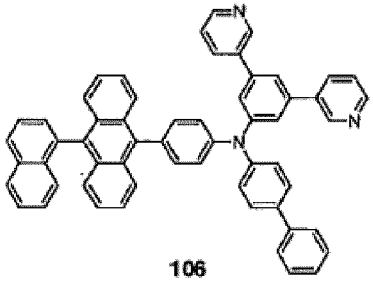

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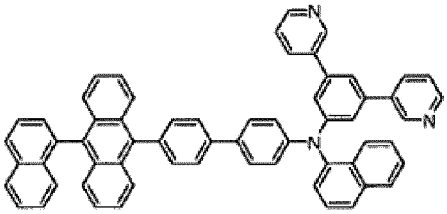

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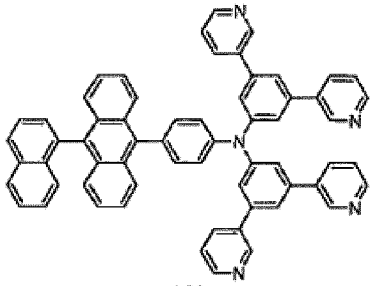

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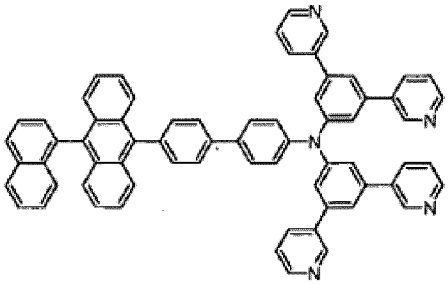


109

19. Organische lichtemittierende Diodenvorrichtung, umfassend eine erste Elektrode, eine gegenüber der ersten Elektrode angeordnete zweite Elektrode und eine zwischen der ersten Elektrode und der zweiten Elektrode angeordnete organische Schicht, wobei die organische Schicht mindestens eine der aminbasierten Verbindungen nach den

Ansprüchen 1 bis 18 umfasst.

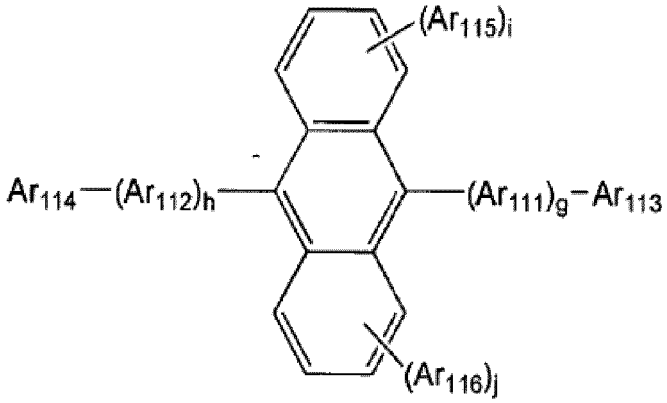
20. Organische lichtemittierende Diodenvorrichtung nach Anspruch 19, wobei die organische Schicht mindestens eine von einer Lochinjektionsschicht, einer Lochtransportschicht, einer funktionellen Schicht mit sowohl Lochinjektions- als auch Lochtransportvermögen, einer Pufferschicht, einer Elektronensperrschicht, einer Emissionsschicht, einer Lochsperrschicht, einer Elektronentransportschicht, einer Elektroneninjektionsschicht und einer funktionellen Schicht mit sowohl Elektroneninjektionsals auch Elektronentransportvermögen umfasst.

21. Organische lichtemittierende Diodenvorrichtung nach Anspruch 20, wobei die organische Schicht eine Elektronentransportschicht umfasst und die aminbasierte Verbindung in der Elektronentransportschicht enthalten ist, vorzugsweise
wobei die Elektronentransportschicht weiterhin einen Metallkomplex umfasst.

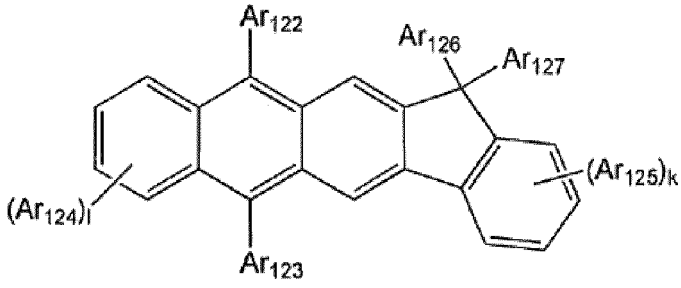
22. Organische lichtemittierende Diodenvorrichtung nach Anspruch 20, wobei die organische Schicht eine Emissionsschicht umfasst und die aminbasierte Verbindung in der Emissionsschicht enthalten ist, vorzugsweise
wobei die aminbasierte Verbindung in der Emissionsschicht als Hostmaterial fungiert und die Emissionsschicht weiterhin einen blau fluoreszierenden Dotanden umfasst.

23. Organische lichtemittierende Diodenvorrichtung nach Anspruch 22, wobei die aminbasierte Verbindung in der Emissionsschicht als Dotand fungiert und die Emissionschicht weiterhin mindestens eine einer durch nachstehende Formel 400 dargestellten Verbindung auf Anthracenbasis und einer durch nachstehende Formel 401 dargestellten Verbindung auf Anthracenbasis umfasst:

Formel 400



Formel 401



wobei, in den Formeln 400 und 401, Ar_{111} und Ar_{112} jeweils unabhängig voneinander ein substituierter oder unsubstituierter C_6-C_{60}-Arylenrest sind; Ar_{113} bis Ar_{116} und Ar_{122} bis Ar_{125} jeweils unabhängig voneinander ein substituierter oder unsubstituierter C_1-C_{10}-Alkylrest oder ein substituierter oder unsubstituierter C_6-C_{60}-Arylrest sind; Ar_{126}

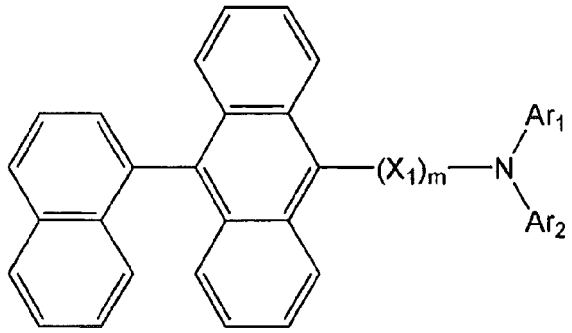
und Ar_{127} jeweils unabhängig voneinander ein C_1-C_{10}-Alkylrest sind; und g, h, i, j, k und l jeweils unabhängig voneinander eine ganze Zahl von 0 bis 4 sind.

24. Organische lichtemittierende Diodenvorrichtung nach Anspruch 20, wobei die organische Schicht mindestens eine von einer Lochinjektionsschicht, einer Lochtransportschicht und einer funktionellen Schicht mit sowohl Lochinjektions- als auch Lochtransportvermögen umfasst und mindestens eine von der Lochinjektionsschicht, der Lochtransportschicht und der funktionellen Schicht mit sowohl Lochinjektions- als auch Lochtransportvermögen eine p-Dotierung umfasst.

Revendications

1. Composé à base d'amine représenté par la Formule 1 ci-dessous :

Formule 1



dans lequel, dans la Formule 1, Ar_1 et Ar_2 sont chacun indépendamment un groupe aryle en C_6 à C_{60} substitué ou non ou un groupe hétéroaryle en C_2 à C_{60} substitué ou non ; et dans lequel Ar_1 et Ar_2 sont éventuellement liés par une liaison simple ;

X_1 est un groupe arylène en C_6 à C_{60} substitué ou non ou un groupe hétéroarylène en C_2 à C_{60} substitué ou non ; m est un nombre entier de 1 à 5 ; et

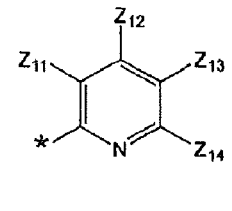
au moins un substituant de chacun du groupe aryle en C_6 à C_{60} substitué, du groupe hétéroaryle en C_2 à C_{60} substitué, du groupe arylène en C_6 à C_{6o} substitué, et du groupe hétéroarylène en C_2 à C_{60} substitué est l'un parmi un atome de deutérium ; -F ; -Cl ; -Br ; -I ; -CN ; un groupe hydroxyle ; $-NO_2$; un groupe amino ; un groupe amidino ; une hydrazine ; une hydrazone ; un groupe carboxyle ou un de ses sels ; un groupe acide sulfonique ou un de ses sels ; un acide phosphorique ou un de ses sels ; un groupe triaryl (C_6 à C_{60}) silyle ; un groupe alkyle en C_1 à C_{60}, un groupe alkoxy en C_1 à C_{60}, un groupe alcényle en C_2 à C_{60}, un groupe alcynyle en C_2 à C_{60} ; un groupe alkyle en C_1 à C_{60}, un groupe alkoxy en C_1 à C_{60}, un groupe alcényle en C_2 à C_{60} et un groupe alcynyle en C_2 à C_{60} qui est substitué par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels ; un groupe cycloalkyle en C_3 à C_{60}, un groupe cycloalcényle en C_3 à C_{60}, un groupe aryle en C_6 à C_{60}, un groupe hétéroaryle en C_2 à C_{60}, un groupe aralkyle en C_6 à C_{60}, un groupe aryloxy en C_6 à C_{60}, un groupe arylthio en C_6 à C_{60} ; un groupe cycloalkyle en C_3 à C_{60}, un groupe cycloalcényle en C_3 à C_{60}, un groupe aryle en C_6 à C_{60}, un groupe hétéroaryle en C_2 à C_{60}, un groupe aralkyle en C_6 à C_{60}, un groupe aryloxy en C_6 à C_{60} et un groupe arylthio en C_6 à C_{60} qui est substitué par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{60}, un groupe alkyle en C_1 à C_{60} substitué par au moins un atome de fluor (F), un groupe alkoxy en C_1 à C_{60}, un groupe alcényle en C_2 à C_{60}, un groupe alcynyle en C_2 à C_{60}, un groupe aryle en C_6 à C_{60} et un groupe hétéroaryle en C_2 à C_{60},

dans lequel au moins un des Ar_1 et Ar_2 est un groupe aryle en C_6 à C_{60} substitué par au moins un groupe attracteur d'électrons choisi dans le groupe constitué par -F ; -CN ; $-NO_2$; un groupe alkyle en C_l à C_{60} substitué par au moins un -F ; un groupe hétéroaryle en C_2 à C_{60} ; et un groupe hétéroaryle en C_2 à C_{60} substitué par un moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino,

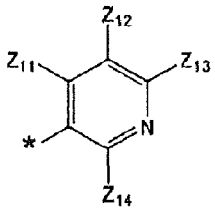
un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels ; un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{60}, un groupe alkyle en C_1 à C_{60} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{60}, un groupe alcényle en C_2 à C_{60}, un groupe alcynyle en C_2 à C_{60}, un groupe aryle en C_6 à C_{60}, et un groupe hétéroaryle en C_2 à C_{60}.

2. Composé à base d'amine selon la revendication 1, dans lequel ledit au moins groupe attracteur d'électrons est choisi dans le groupe constitué par : -F ; -CN ; $-NO_2$; un groupe alkyle en C_1 à C_{20} substitué par au moins un -F ; un groupe hétéroaryle en C_2 à C_{20} comprenant un noyau contenant un atome N ; et un groupe hétéroaryle en C_2 à C_{20} qui comprend un noyau contenant un atome N et est substitué par un moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{20}, un groupe alkyle en C_1 à C_{20} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{20}.
un groupe aryle en C_6 à C_{20}, et un groupe hétéroaryle en C_2 à C_2.

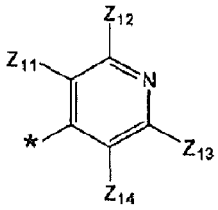
3. Composé à base d'amine selon la revendication 1 ou la revendication 2, dans lequel ledit au moins un groupe attracteur d'électrons est choisi dans le groupe constitué par -F ; -CN ; $-CH_2F$; $-CHF_2$; $-CF_3$; et les groupes représentés par les Formules 2(1) à 2(14) ci-dessous :



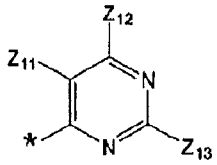
Formule 2(1)



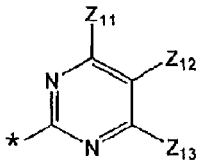
Formule 2(2)



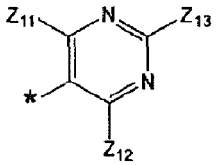
Formule 2(3)



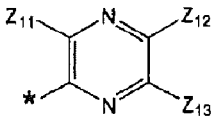
Formule 2(4)



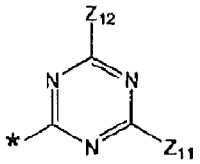
Formule 2(5)



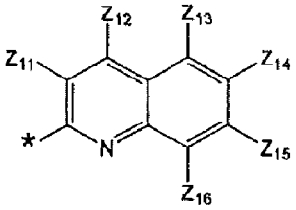
Formule 2(6)



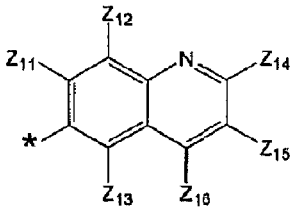
Formule 2(7)



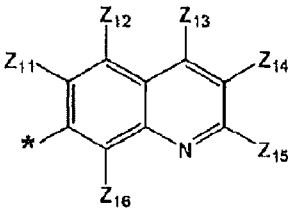
Formule 2(8)



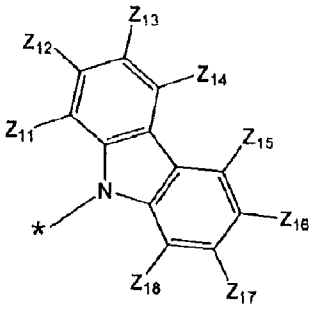
Formule 2(9)



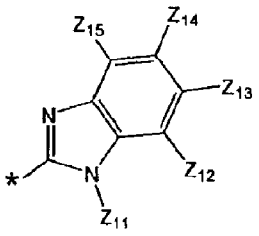
Formule 2(10)



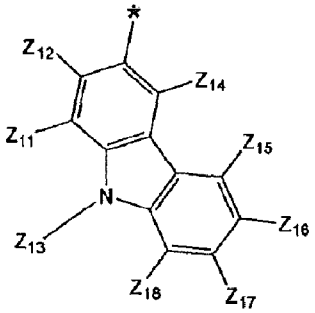
Formule 2(11)



Formule 2(12)



Formule 2(13)



Formule 2(14)

dans lequel, dans les Formules 2(1) à 2(14), Z_{11} à Z_{18} sont chacun indépendamment un atome d'hydrogène, un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{20}, un groupe alkyle en C_1 à C_{20} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{20}, un groupe phényle, un groupe naphtyle, un groupe anthryle, un groupe phénanthrényle, un groupe pyrényle, un groupe pyridinyle, un groupe triazinyle, ou un groupe carbazolyle.

4. Composé à base d'amine selon l'une quelconque des revendications précédentes, dans lequel Ar_1 et Ar_2 sont chacun indépendamment un groupe aryle en C_6 à C_{60} choisi parmi un groupe phényle substitué ou non, un groupe naphtyle substitué ou non, un groupe fluorényle substitué ou non, un groupe phénanthrényle substitué ou non, un groupe anthryle substitué ou non, un groupe triphénylényle substitué ou non, un groupe pyrényle substitué ou non, un groupe chrysényle substitué ou non, un groupe pyridinyle substitué ou non, un groupe pyrazinyle substitué ou non, un groupe pyrimidinyle substitué ou non, un groupe quinolyle substitué ou non, un groupe carbazolyle substitué ou non, un groupe triazinyle substitué ou non, un groupe dibenzothiophényle substitué ou non, un groupe dibenzofurannyle substitué ou non, ou un groupe phénanthrolinyle substitué ou non.

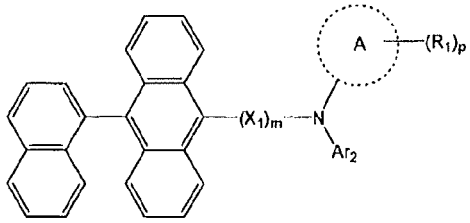
5. Composé à base d'amine selon l'une quelconque des revendications précédentes, dans lequel Ar_1 et Ar_2 sont liés par une liaison simple.

6. Composé à base d'amine selon l'une quelconque des revendications précédentes, dans lequel ledit au moins Ar_1 et Ar_2 est un groupe aryle en C_6 à C_{60} substitué par au moins deux groupes attracteurs d'électrons.

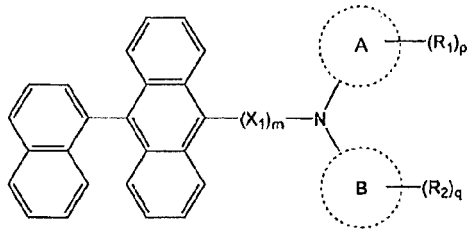
7. Composé à base d'amine selon l'une quelconque des revendications précédentes, dans lequel ledit au moins Ar_1 et Ar_2 est un groupe phényle, un groupe biphényle, un groupe naphtyle, un groupe anthryle, un groupe phénanthrényle, un groupe pyrényle, ou un groupe fluorényle qui est substitué par au moins deux groupes attracteurs d'électrons ; et
les groupes attracteurs d'électrons sont chacun indépendamment choisis dans le groupe constitué par un groupe pyridinyle, un groupe pyrazinyle, un groupe pyrimidinyle, un groupe quinolinyle, un groupe isoquinolinyle, un groupe quinazolinyle, un groupe triazinyle, un groupe benzoimidazolyle, et un groupe carbazolyle ; et un groupe pyridinyle, un groupe pyrazinyle, un groupe pyrimidinyle, un groupe quinolinyle, un groupe isoquino-linyle, un groupe quinazolinyle, un groupe phtalazinyle, un groupe benzoimidazolyle, et un groupe carbazolyle qui sont substitués par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN ; un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{20}, un groupe alkyle en C_1 à C_{20} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{20}, un groupe phényle, un groupe naphtyle, un groupe anthryle, un groupe phénanthrényle, un groupe pyrényle, un groupe pyridinyle, un groupe triazinyle, et un groupe carbazolyle.

8. Composé à base d'amine selon la revendication 1, dans lequel le composé à base d'amine est représenté par la Formule 1(1) ou 1(2) ci-dessous :

Formule 1(1)



Formule 1(2)



dans lequel, dans la Formule 1(1), Ar_2 est un groupe aryle en C_6 à C_{20} substitué ou non ou un groupe hétéroaryle en C_2 à C_{20} substitué ou non ; et
dans les Formules 1(1) et 1(2), les groupes à noyau A et B sont chacun indépendamment un groupe aryle en C_6 à C_{20} substitué ;
R_1 et R_2 sont chacun indépendamment un groupe attracteur d'électrons choisi dans le groupe constitué par -F ; -CN ; $-NO_2$; un groupe alkyle en C_1 à C_{60} substitué par au moins un -F ; un groupe hétéroaryle en C_2 à C_{60} ; et un groupe hétéroaryle en C_2 à C_{60} substitué par un moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{60}, un groupe alkyle en C_1 à C_{60} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{60}, un groupe alcényle en C_2 à C_{60}, un groupe alcynyle en C_2 à C_{60}, un groupe aryle en C_6 à C_{60}, et un groupe hétéroaryle en C_2 à C_{60} ; et
p et q sont chacun indépendamment un nombre entier de 1 à 9.

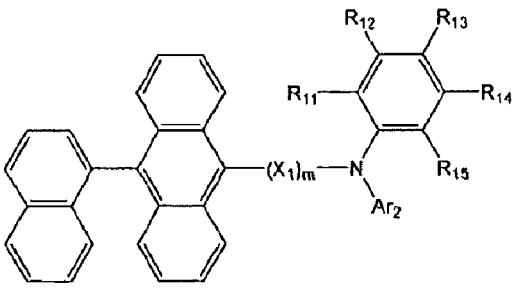
9. Composé à base d'amine selon la revendication 8, dans lequel le composé à base d'amine est représenté par la Formule 1(1) dans laquelle au moins un de R_1 est -CN ; ou est représenté par la Formule 1(2) dans laquelle au moins un de R_1 et un de R_2 est -CN.

10. Composé à base d'amine selon la revendication 7, dans lequel le composé à base d'amine est représenté par la Formule 1(1), dans laquelle le groupe à noyau A est un groupe phényle substitué, un groupe biphényle substitué, un groupe naphtyle substitué, un groupe anthryle substitué, un groupe phénanthrényle substitué, un groupe pyrényle substitué, ou un groupe florényle substitué ;
R_1 est au moins un groupe attracteur d'électrons choisi dans le groupe constitué par un groupe pyridinyle, un groupe pyrazinyle, un groupe pyrimidinyle, un groupe quinolinyle, un groupe isoquinolinyle, un groupe quinazolinyle, un groupe triazinyle, un groupe benzoimidazolyle, et un groupe carbazolyle ; et un groupe pyridinyle, un groupe pyrazinyle, un groupe pyrimidinyle, un groupe quinolinyle, un groupe isoquinolinyle, un groupe quinazolinyle, un groupe phtalazinyle, un groupe benzoimidazolyle, et un groupe carbazolyle qui sont substitués par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN ; un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{20}, un groupe alkyle en C_1 à C_{20} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{20}, un groupe phényle, un groupe naphtyle, un groupe anthryle, un groupe phénanthrényle, un groupe pyrényle, un groupe pyridinyle, un groupe triazinyle, et un groupe carbazolyle ; et p est 2, 3, ou 4 ; ou, dans lequel le composé à base d'amine est représenté par la Formule 1(2), dans laquelle le groupe à noyau A et le groupe à noyau B sont chacun indépendamment un groupe phényle substitué, un groupe biphényle substitué, un groupe naphtyle substitué, un groupe anthryle substitué, un groupe phénanthrényle substitué, un groupe pyrényle substitué, ou un groupe florényle substitué ;
R_1 et R_2 sont chacun indépendamment au moins un groupe attracteur d'électrons choisi dans le groupe constitué

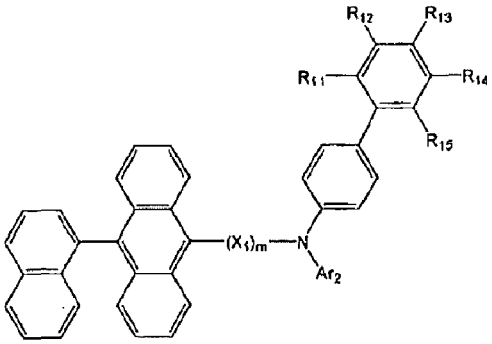
par un groupe pyridinyle, un groupe pyrazinyle, un groupe pyrimidinyle, un groupe quinolinyle, un groupe isoquinolinyle, un groupe quinazolinyle, un groupe triazinyle, un groupe benzoimidazolyle, et un groupe carbazolyle ; et un groupe pyridinyle, un groupe pyrazinyle, un groupe pyrimidinyle, un groupe quinolinyle, un groupe isoquinolinyle, un groupe quinazolinyle, un groupe phtalazinyle, un groupe benzoimidazolyle, et un groupe carbazolyle qui sont substitués par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN ; un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{20}, un groupe alkyle en C_1 à C_{20} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{20}, un groupe phényle, un groupe naphtyle, un groupe anthryle, un groupe phénanthrényle, un groupe pyrényle, un groupe pyridinyle, un groupe triazinyle, et un groupe carbazolyle ; et p et q sont chacun indépendamment 2, 3, ou 4.

11. Composé à base d'amine selon la revendication 1, dans lequel le composé à base d'amine est représenté par une des Formules 1A à 1J ci-dessous :

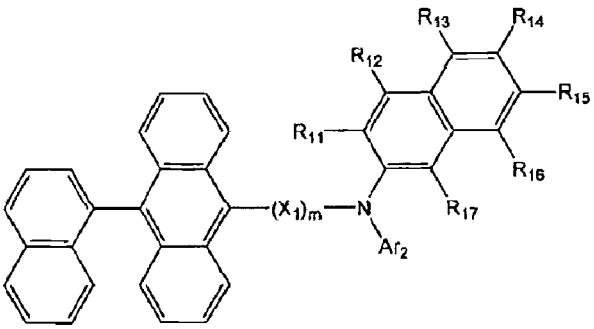
Formule 1A



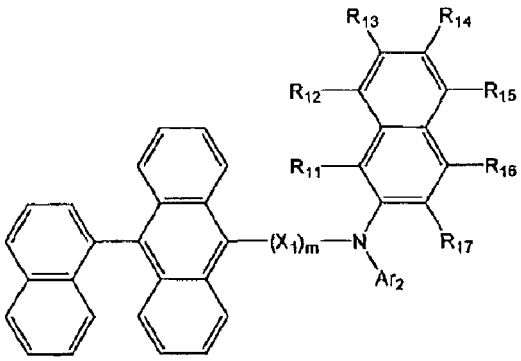
Formule 1B



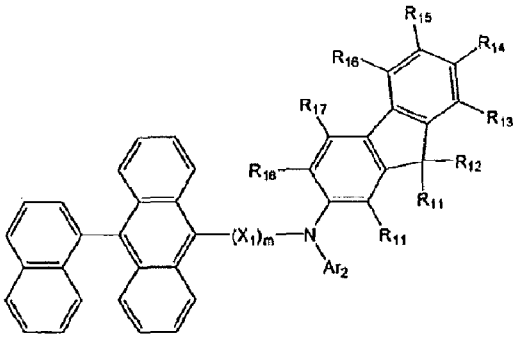
Formule 1C



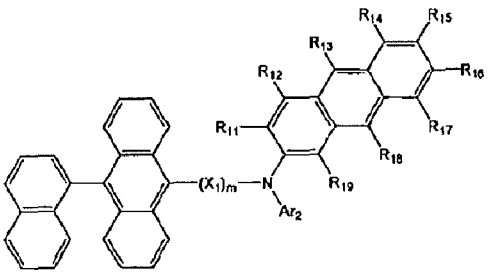
Formule 1D



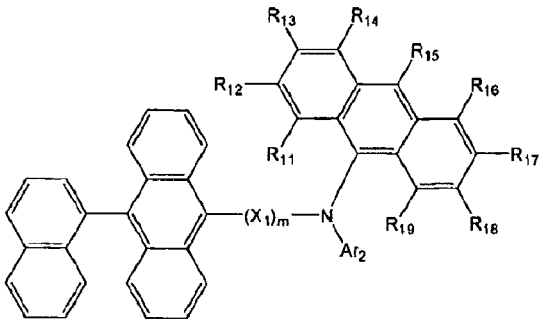
Formule 1E



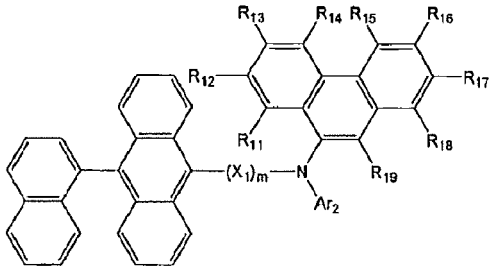
Formule 1F



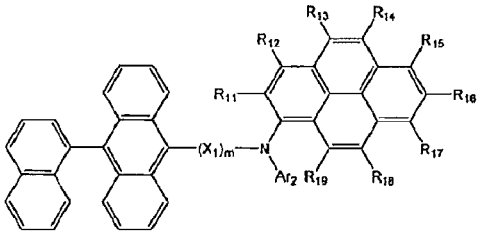
Formule 1G



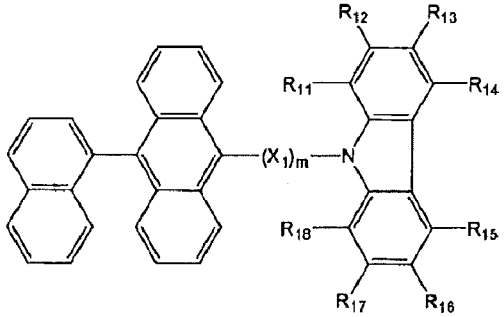
Formule 1H



Formule 1I



Formule 1J



dans lequel, dans les Formules 1A à 1J, Ar_2 est un groupe phényle substitué ou non, un groupe pentalényle substitué ou non, un groupe indénylène substitué ou non, un groupe naphtyle substitué ou non, un groupe azulényle substitué ou non, un groupe heptalényle substitué ou non, un groupe indacényle substitué ou non, un groupe acénaphtyle substitué ou non, un groupe fluorényle substitué ou non, un groupe phénalényle substitué ou non, un groupe phénanthrényle substitué ou non, un groupe anthryle substitué ou non, un groupe fluoranthényle substitué ou non, un groupe triphénylényle substitué ou non, un groupe pyrényle substitué ou non, un groupe chrysényle substitué ou non, un groupe naphtacényle substitué ou non, un groupe picényle substitué ou non, un groupe pérylényle substitué ou non, un groupe pentaphényle substitué ou non, un groupe hexacényle substitué ou non, un groupe pyrrolyle substitué ou non, un groupe pyrazolyle substitué ou non, un groupe imidazolyle substitué ou non, un groupe imidazolinyle substitué ou non, un groupe imidazopyridinyle substitué ou non, un groupe imidazopyrimidinyle substitué ou non, un groupe pyridinyle substitué ou non, un groupe pyrazinyle substitué ou non, un groupe pyrimidinyle substitué ou non, un groupe benzoimidazolyle substitué ou non, un groupe indolyle substitué ou non, un groupe purinyle substitué ou non, un groupe quinolinyle substitué ou non, un groupe phtalazinyle substitué ou non, un groupe indolizinyle substitué ou non, un groupe naphtyridinyle substitué ou non, un groupe quinazolinyle substitué ou non, un groupe cinolinyle substitué ou non, un groupe indazolyle substitué ou non, un groupe carbazolyle substitué ou non, un groupe phénazinyle substitué ou non, un groupe phénanthridinyle substitué ou non, un groupe pyranyle substitué ou non, un groupe chroménylе substitué ou non, un groupe furannyle substitué ou non, un groupe benzofurannyle substitué ou non, un groupe thiophényle substitué ou non, un groupe benzothiophényle substitué ou non, un groupe isothiazolyle substitué ou non, un groupe benzoimidazolyle substitué ou non, un groupe isoxazolyle substitué ou non, un groupe dibenzothiophényle substitué ou non, un groupe dibenzofurannyle substitué ou non, un groupe triazinyle substitué ou non, un groupe oxadiazolyle substitué ou non, un groupe pyridazinyle substitué ou non, un groupe triazolyle substitué ou non, un groupe tétrazolyle substitué ou non, ou un groupe phénanthrolinyle substitué ou non,

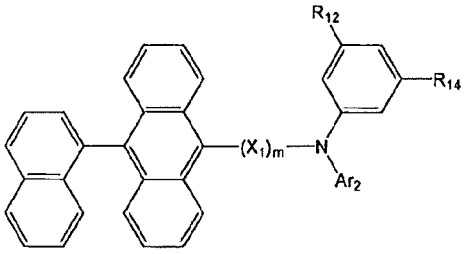
les substituants du groupe phényle substitué, du groupe pentalényle substitué, du groupe indényle substitué, du groupe naphtyle substitué, du groupe azulényle substitué, du groupe heptalényle substitué, du groupe indacényle substitué, du groupe acénaphtyle substitué, du groupe fluorényle substitué, du groupe phénalényle substitué, du groupe phénanthrényle substitué, du groupe anthryle substitué, du groupe fluoranthényle substitué, du groupe triphénylényle substitué, du groupe pyrényle substitué, du groupe chrysényle substitué, du groupe naphtacényle substitué, du groupe picényle substitué, du groupe pérylényle substitué, du groupe pentaphényle substitué, du groupe hexacényle substitué, du groupe pyrrolyle substitué, du groupe pyrazolyle substitué, du groupe imidazolyle substitué, du groupe imidazolinyle substitué, du groupe imidazopyridinyle substitué, du groupe imidazopyrimidinyle

substitué, du groupe pyridinyle substitué, du groupe pyrazinyle substitué, du groupe pyrimidinyle substitué, du groupe benzoimidazolyle substitué, du groupe indolyle substitué, du groupe purinyle substitué, du groupe quinolinyle substitué, du groupe phtalazinyle substitué, du groupe indolizinyle substitué, du groupe naphtyridinyle substitué, du groupe quinazolinyle substitué, du groupe cinolinyle substitué, du groupe indazolyle substitué, du groupe carbazolyle substitué, du groupe phénazinyle substitué, du groupe phénanthridinyle substitué, du groupe pyranyle substitué, du groupe chroménylе substitué, du groupe furannyle substitué, du groupe benzofurannyle substitué, du groupe thiophényle substitué, du groupe benzothiophényle substitué, du groupe isothiazolyle substitué, du groupe benzoimidazolyle substitué, du groupe isoxazolyle substitué, du groupe dibenzothiophényle substitué, du groupe dibenzofurannyle substitué, du groupe triazinyle substitué, du groupe oxadiazolyle substitué, du groupe pyridazinyle substitué, du groupe triazolyle substitué, du groupe tétrazolyle substitué, et du groupe phénanthrolinyle substitué, et R_{11} à R_{19} sont chacun indépendamment un atome d'hydrogène ; un atome de deutérium ; -F ; -Cl ; -Br ; -I ; -CN ; un groupe hydroxyle ; $-NO_2$; un groupe amino ; un groupe amidino ; une hydrazine ; une hydrazone ; un groupe carboxyle ou un de ses sels ; un groupe acide sulfonique ou un de ses sels ; un acide phosphorique ou un de ses sels ; un groupe triaryl (C_6 à C_{60}) silyle ; un groupe alkyle en C_1 à C_{60}, un groupe alkoxy en C_1 à C_{60}, un groupe alcényle en C_2 à C_{60}, et un groupe alcynyle en C_2 à C_{60} ; un groupe alkyle en C_1 à C_{60}, un groupe alkoxy en C_1 à C_{60}, un groupe alcényle en C_2 à C_{60} et un groupe alcynyle en C_2 à C_{60} qui sont substitués par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, et un acide phosphorique ou un de ses sels ; un groupe cycloalkyle en C_3 à C_{60}, un groupe cycloalcényle en C_3 à C_{60}, un groupe aryle en C_6 à C_{60}, un groupe hétéroaryle en C_2 à C_{60}, un groupe aralkyle en C_6 à C_{60}, un groupe aryloxy en C_6 à C_{60}, un groupe arylthio en C_6 à C_{60} ; un groupe cycloalkyle en C_3 à C_{60}, un groupe cycloalcényle en C_3 à C_{60}, un groupe aryle en C_6 à C_{60}, un groupe hétéroaryle en C_2 à C_{60}, un groupe aralkyle en C_6 à C_{60}, un groupe aryloxy en C_6 à C_{60}, et un groupe arylthio en C_6 à C_{60} qui sont substitués par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{60}, un groupe alkyle en C_1 à C_{60} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{60}, un groupe alcényle en C_2 à C_{60}, un groupe alcynyle en C_2 à C_{60}, un groupe aryle en C_6 à C_{60} et un groupe hétéroaryle en C_2 à C_{60},

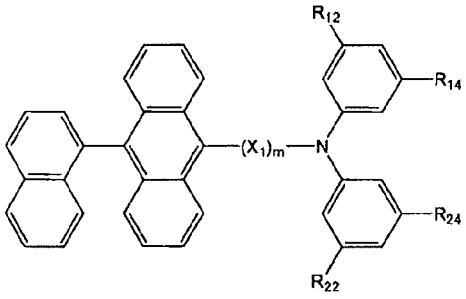
dans lequel au moins un des R_{11} à R_{15} dans les Formules 1A et 1B, au moins un des R_{11} à R_{17} dans les Formules 1C et 1D, au moins un des R_{11} à R_{18} dans les Formules 1E et 1J, et au moins un des R_{11} à R_{19} dans la Formule 1F, 1G, 1H et 1I est indépendamment un groupe attracteur d'électrons choisi dans le groupe constitué par -F ; -CN ; $-NO_2$; un groupe alkyle en C_1 à C_{60} substitué par au moins un -F ; un groupe hétéroaryle en C_2 à C_{60} ; et un groupe hétéroaryle en C_2 à C_{60} substitué par un moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{60}, un groupe alkyle en C_1 à C_{60} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{60}, un groupe alcényle en C_2 à C_{60}, un groupe alcynyle en C_2 à C_{60}, un groupe aryle en C_6 à C_{60}, et un groupe hétéroaryle en C_2 à C_{60}.

12. Composé à base d'amine selon la revendication 11, dans lequel le composé à base d'amine est représenté par la Formule 1A-(1) ou 1A-(2) ci-dessous :

Formule 1A-(1)



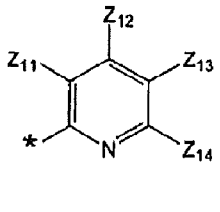
Formule 1A-(2)



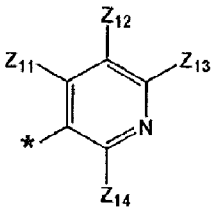
dans lequel, dans les Formules 1A-(1) et 1A-(2), R_{12}, R_{14}, R_{22}, et R_{24} sont chacun indépendamment un groupe attracteur d'électrons choisi dans le groupe constitué par un groupe pyridinyle, un groupe pyrazinyle, un groupe pyrimidinyle, un groupe quinolinyle, un groupe isoquinolinyle, un groupe quinazolinyle, un groupe triazinyle, un

groupe benzoimidazolyle, et un groupe carbazolyle ; et un groupe pyridinyle, un groupe pyrazinyle, un groupe pyrimidinyle, un groupe quinolinyle, un groupe isoquinolinyle, un groupe quinazolinyle, un groupe phtalazinyle, un groupe benzoimidazolyle, et un groupe carbazolyle qui sont substitués par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN ; un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{20}, un groupe alkyle en C_1 à C_{20} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{20}, un groupe phényle, un groupe naphtyle, un groupe anthryle, un groupe phénanthrényle, un groupe pyrényle, un groupe pyridinyle, un groupe triazinyle, et un groupe carbazolyle ; et
dans la Formule 1A-(1), Ar_2 est un groupe phényle substitué ou non, un groupe biphényle substitué ou non, un groupe naphtyle substitué ou non, un groupe anthryle substitué ou non, un groupe phénanthrényle substitué ou non, un groupe pyrényle substitué ou non, ou un groupe fluorényle substitué ou non.

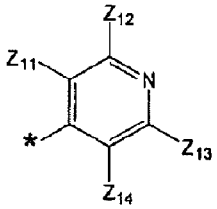
13. Composé à base d'amine selon la revendication 12, dans lequel R_{12}, R_{14}, R_{22} et R_{24} sont chacun indépendamment choisis dans le groupe constitué par -F ; -CN ; $-CH_2F$; $-CHF_2$; $-CF_3$; et les groupes représentés par les Formules 2(1) à 2(14) ci-dessous :



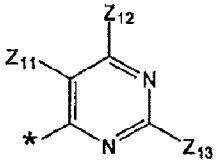
Formule 2(1)



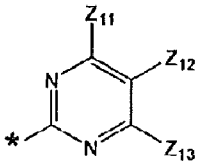
Formule 2(2)



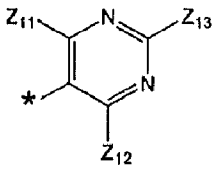
Formule 2(3)



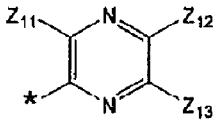
Formule 2(4)



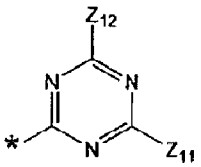
Formule 2(5)



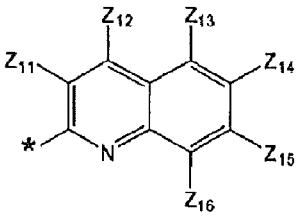
Formule 2(6)



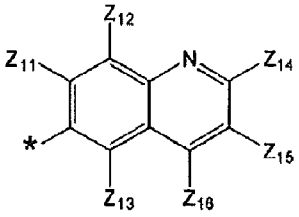
Formule 2(7)



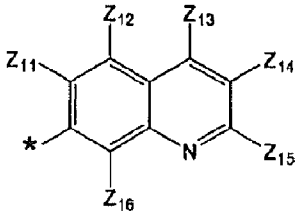
Formule 2(8)



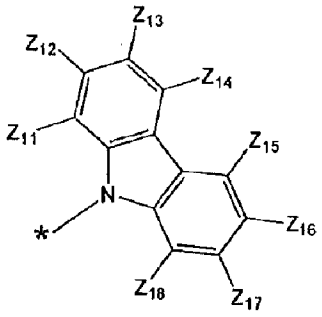
Formule 2(9)



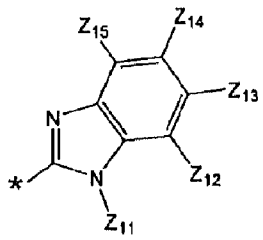
Formule 2(10)



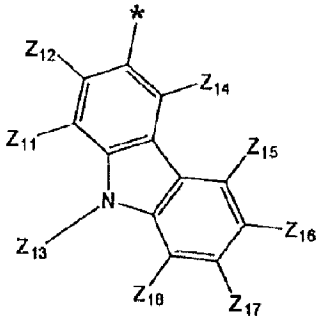
Formule 2(11)



Formule 2(12)



Formule 2(13)

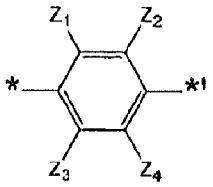


Formule 2(14)

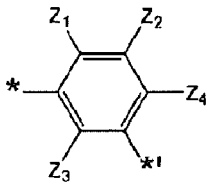
dans lequel, dans les Formules 2(1) à 2(14), Z_{11} à Z_{18} sont chacun indépendamment un atome d'hydrogène, un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{20}, un groupe alkyle en C_1 à C_{20} substitué par au moins un -F, un groupe alkoxy en C_1 à C_{20}, un groupe phényle, un groupe naphtyle, un groupe anthryle, un groupe phénanthrényle, un groupe pyrényle, un groupe pyridinyle, un groupe triazinyle, ou un groupe carbazolyle.

14. Composé à base d'amine selon l'une quelconque des revendications précédentes, dans lequel X_1 est un groupe phénylène substitué ou non, un groupe pentalénylène substitué ou non, un groupe indénylène substitué ou non, un groupe naphtylène substitué ou non, un groupe azulénylène substitué ou non, un groupe heptalénylène substitué ou non, un groupe indacénylène substitué ou non, un groupe acénaphtylène substitué ou non, un groupe fluorénylène substitué ou non, un groupe phénalénylène substitué ou non, un groupe phénanthrénylène substitué ou non, un groupe anthrylène substitué ou non, un groupe fluoranthénylène substitué ou non, un groupe triphénylénylène substitué ou non, un groupe pyrénylène substitué ou non, un groupe chrysénylène substitué ou non, un groupe naphtacénylène substitué ou non, un groupe picénylène substitué ou non, un groupe pérylénylène substitué ou non, un groupe pentaphénylène substitué ou non, un groupe hexacénylène substitué ou non, un groupe pyrrolylène substitué ou non, un groupe pyrazolylène substitué ou non, un groupe imidazolylène substitué ou non, un groupe imidazolinylène substitué ou non, un groupe imidazopyridinylène substitué ou non, un groupe imidazopyrimidinylène substitué ou non, un groupe pyridinylène substitué ou non, un groupe pyrazinylène substitué ou non, un groupe pyrimidinylène substitué ou non, un groupe indolylène substitué ou non, un groupe purinylène substitué ou non, un groupe quinolinylène substitué ou non, un groupe phtalazinylène substitué ou non, un groupe indolizinylène substitué ou non, un groupe naphtyridinylène substitué ou non, un groupe quinazolinylène substitué ou non, un groupe cinnolinylène substitué ou non, un groupe indazolylène substitué ou non, un groupe carbazolylène substitué ou non, un groupe phénazinylène substitué ou non, un groupe phénanthridinylène substitué ou non, un groupe pyranylène substitué ou non, un groupe chroménylène substitué ou non, un groupe furanylène substitué ou non, un groupe benzofuranylène substitué ou non, un groupe thiophénylène substitué ou non, un groupe benzothiophénylène substitué ou non, un groupe isothiazolylène substitué ou non, un groupe benzoimidazolylène substitué ou non, un groupe isoxazolylène substitué ou non, un groupe dibenzothiophénylène substitué ou non, un groupe dibenzofuranylène substitué ou non, un groupe triazinylène substitué ou non, un groupe oxadiazolylène substitué ou non, un groupe pyridazinylène substitué ou non, un groupe triazolylène substitué ou non, ou un groupe tétrazolylène substitué ou non.

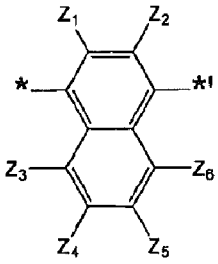
15. Composé à base d'amine selon la revendication 14, dans lequel X_1 est un groupe représenté par une des Formules 5(1) à 5(16) ci-dessous :



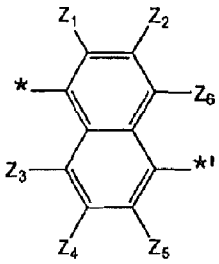
Formule 5(1)



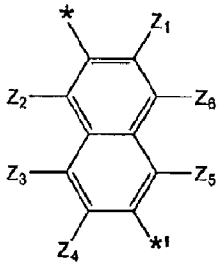
Formule 5(2)



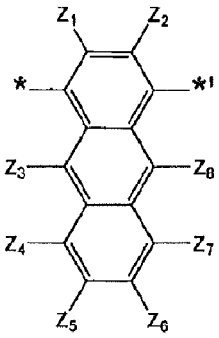
Formule 5(3)



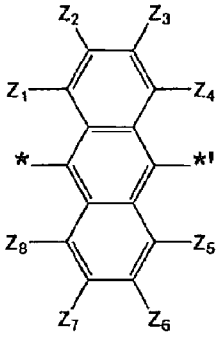
Formule 5(4)



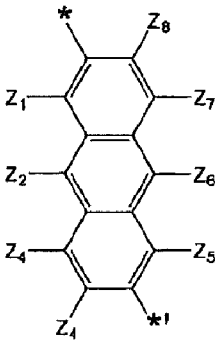
Formule 5(5)



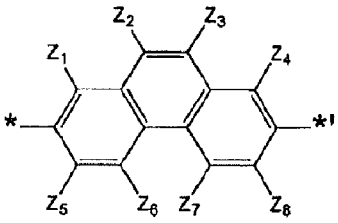
Formule 5(6)



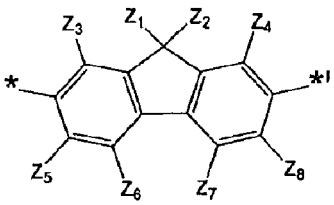
Formule 5(7)



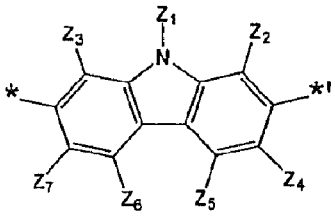
Formule 5(8)



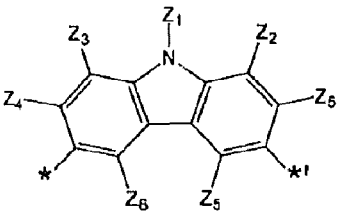
Formule 5(9)



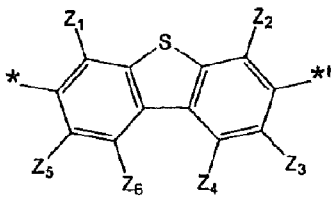
Formule 5(10)



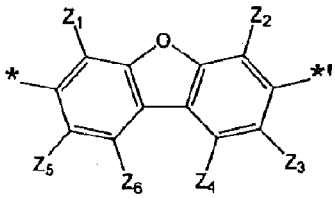
Formule 5(11)



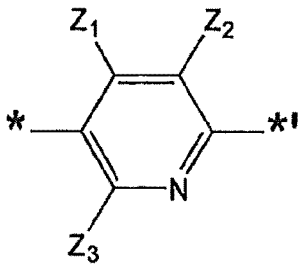
Formule 5(12)



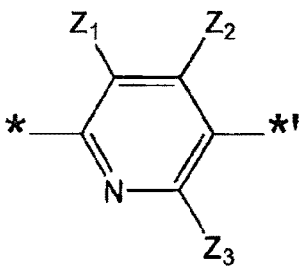
Formule 5(13)



Formule 5(14)



Formule 5(15)



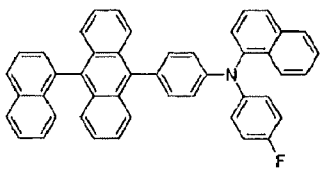
Formule 5(16)

dans lequel, dans les Formules 5(1) à 5(16), Z_1 à Z_8 sont chacun indépendamment l'un parmi un atome d'hydrogène ; un atome de deutérium ; -F ; -Cl ; -Br ; -I ; -CN ; un groupe hydroxyle ; $-NO_2$; un groupe amino ; un groupe amidino ; une hydrazine ; une hydrazone ; un groupe carboxyle ou un de ses sels ; un groupe acide sulfonique ou un de ses sels ; un acide phosphorique ou un de ses sels ; un groupe alkyle en C_1 à C_{20} ; un groupe alkoxy en C_1 à C_{20} ; un groupe alkyle en C_1 à C_{20} et un groupe alkoxy en C_1 à C_{20} qui sont substitués par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, et un acide phosphorique ou un de ses sels ; un groupe aryle en C_6 à C_{20} ; un groupe hétéroaryle en C_2 à C_{20} ; et un groupe aryle en C_6 à C_{20} et un groupe hétéroaryle en C_2 à C_{20} qui sont substitués par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{20}, un groupe alkoxy en C_1 à C_{20}, un groupe aryle en C_6 à C_{60} et un groupe hétéroaryle en C_2 à C_{20} ;
dans lequel * indique un site de liaison à l'anthracène dans la Formule 1 ; et
dans lequel *' indique un site de liaison à N dans la Formule 1.

16. Composé à base d'amine selon la revendication 15, dans lequel Z_1 à Z_8 sont chacun indépendamment l'un parmi un atome d'hydrogène ; un atome de deutérium ; -F ; -Cl ; -Br ; -I ; -CN ; un groupe hydroxyle ; $-NO_2$; un groupe amino ; un groupe amidino ; une hydrazine ; une hydrazone ; un groupe carboxyle ou un de ses sels ; un groupe acide sulfonique ou un de ses sels ; un acide phosphorique ou un de ses sels ; un groupe méthyle, un groupe éthyle, un groupe propyle, un groupe butyle, et un groupe pentyle ; un groupe méthoxy, un groupe éthoxy, un groupe propoxy, un groupe butoxy, et un groupe pentoxy ; un groupe méthyle, un groupe éthyle, un groupe propyle, un groupe butyle, et un groupe pentyle, un groupe méthoxy, un groupe éthoxy, un groupe propoxy, un groupe butoxy, et un groupe pentoxy qui sont substitués par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, et un acide phosphorique ou un de ses sels ; un groupe phényle, un groupe naphtyle, un groupe anthryle, un groupe phénanthrényle, un groupe pyrényle, et un groupe fluorényle ; un groupe pyridinyle, un groupe pyrimidinyle, un groupe triazinyle, un groupe quinolyle, et un groupe carbazolyle ; un groupe phényle, un groupe naphtyle, un groupe anthryle, un groupe phénanthrényle, un groupe pyrényle, et un groupe fluorényle ; un groupe pyridinyle, un groupe pyrimidinyle, un groupe triazinyle, un groupe quinolyle, et un groupe carbazolyle qui sont substitués par au moins l'un parmi un atome de deutérium, -F, -Cl, -Br, -I, -CN, un groupe hydroxyle, $-NO_2$, un groupe amino, un groupe amidino, une hydrazine, une hydrazone, un groupe carboxyle ou un de ses sels, un groupe acide sulfonique ou un de ses sels, un acide phosphorique ou un de ses sels, un groupe alkyle en C_1 à C_{20}, et un groupe alkoxy en C_1 à C_{20}.

17. Composé à base d'amine selon l'une quelconque des revendications précédentes, dans lequel m est 1, 2, ou 3.

18. Composé à base d'amine selon l'une quelconque des revendications précédentes, dans lequel le composé à base d'amine est un des composés 1 à 109 ci-dessous :



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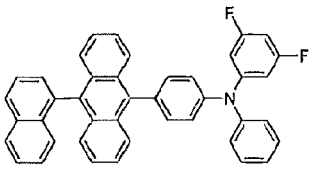

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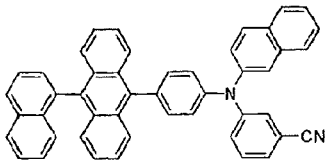

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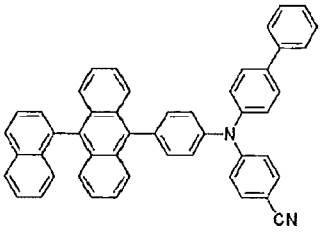

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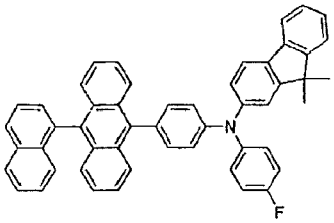

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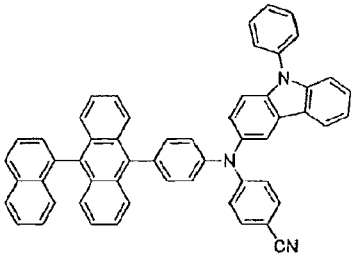

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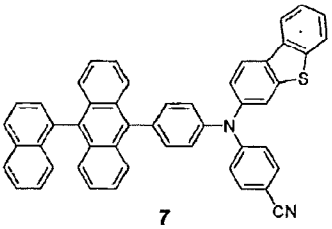

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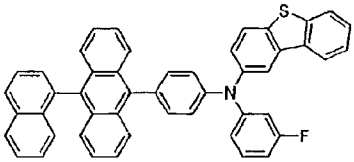

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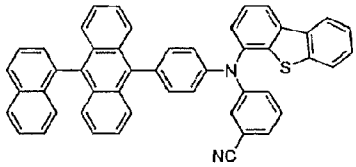

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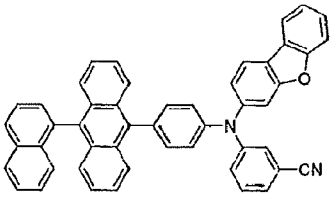

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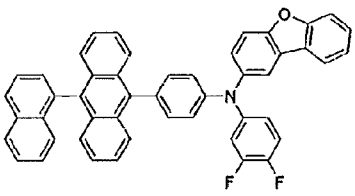

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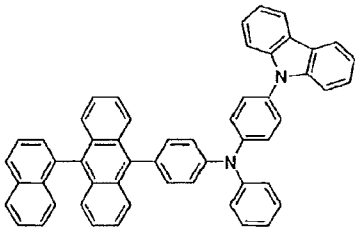

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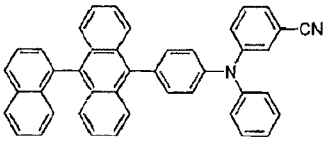

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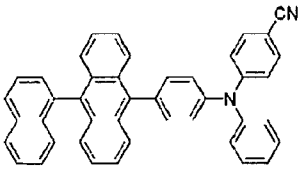

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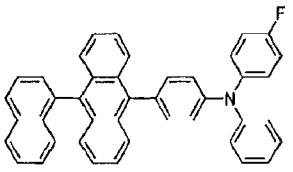

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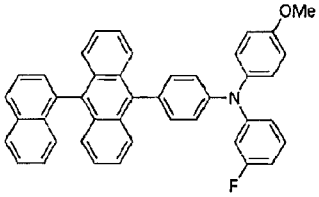

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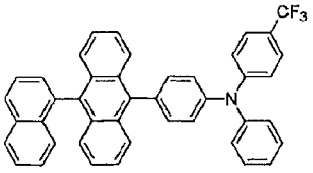

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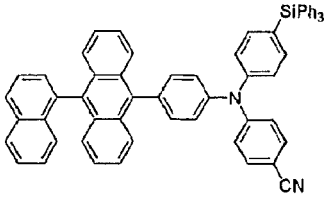

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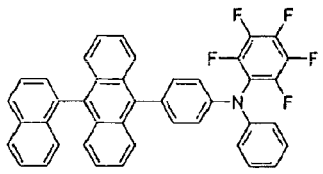

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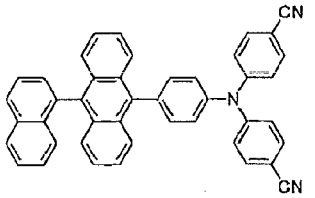

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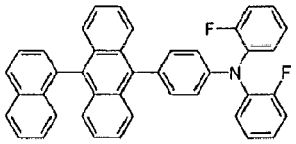

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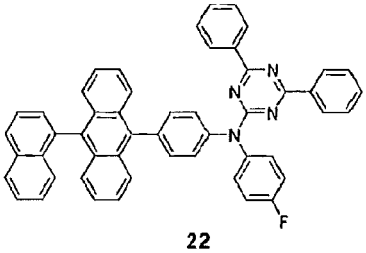

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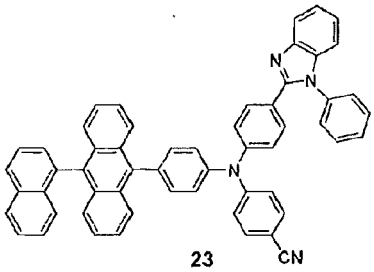

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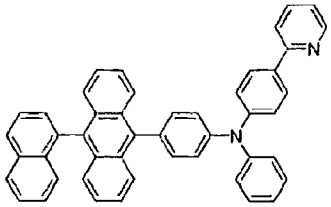

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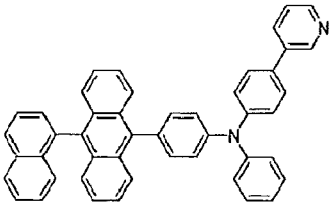

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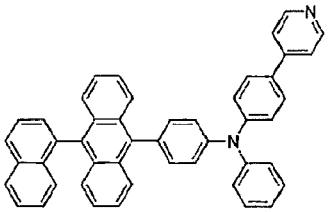

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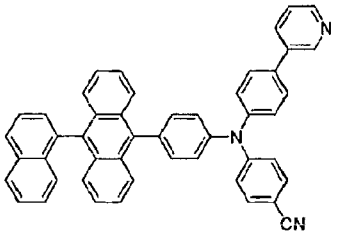

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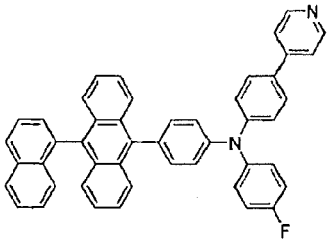

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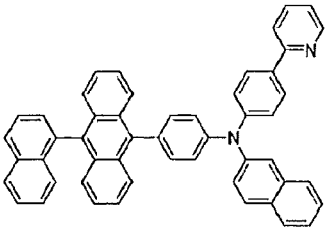

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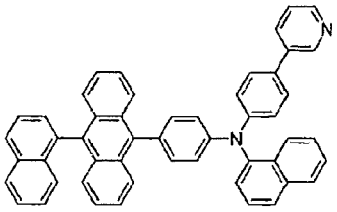

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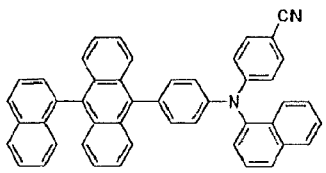

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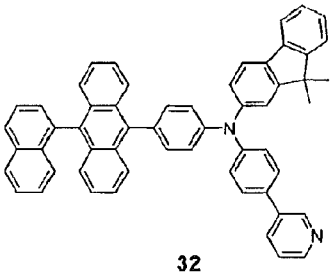

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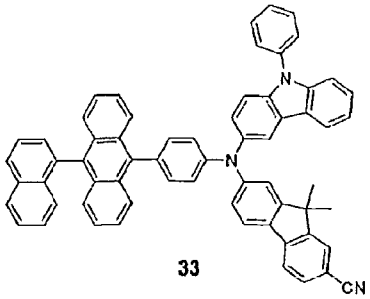

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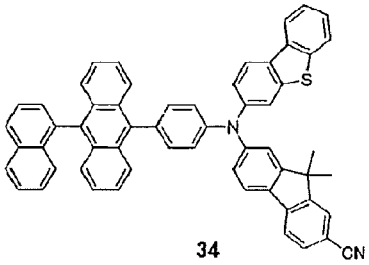

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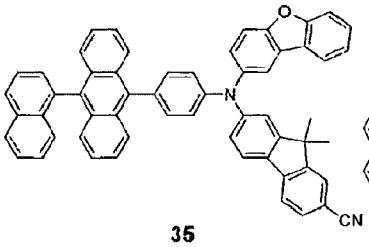

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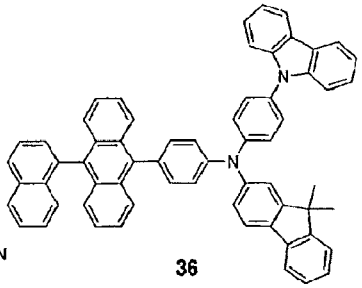

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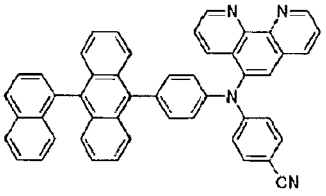

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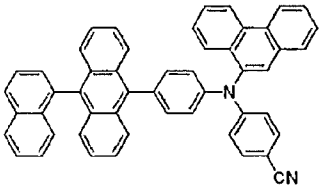

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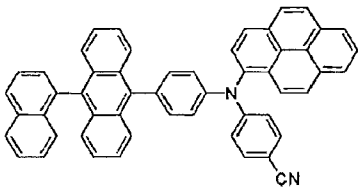

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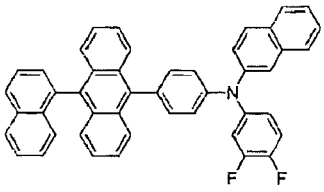

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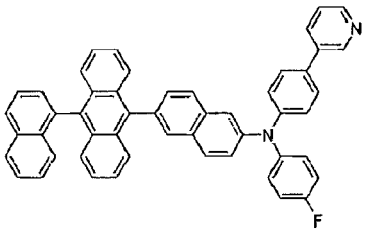

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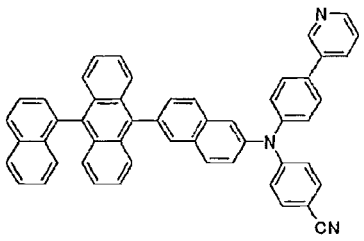

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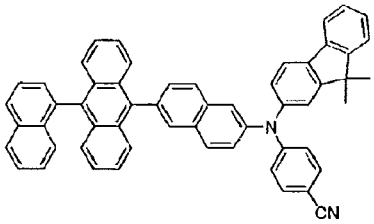

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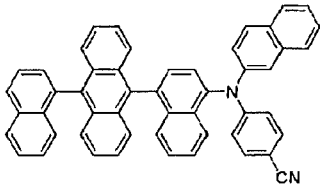

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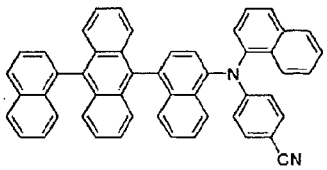

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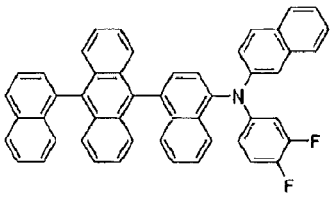

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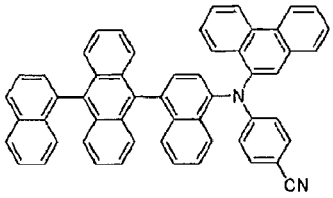

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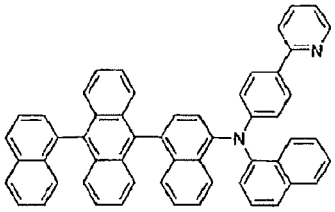

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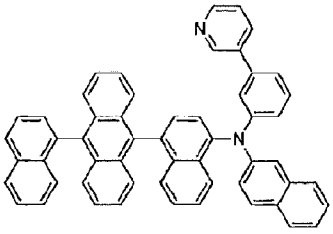

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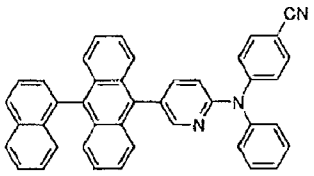

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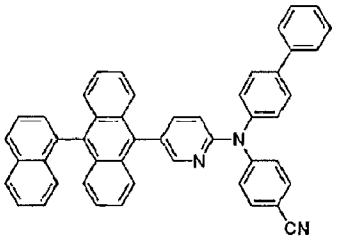

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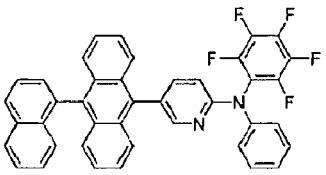

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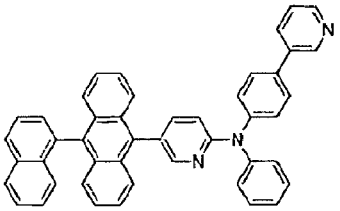

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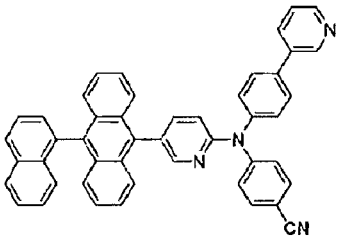

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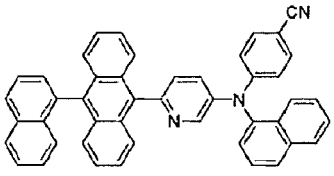

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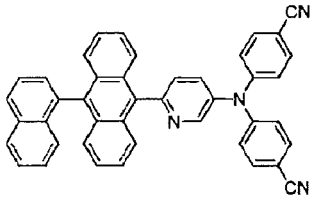

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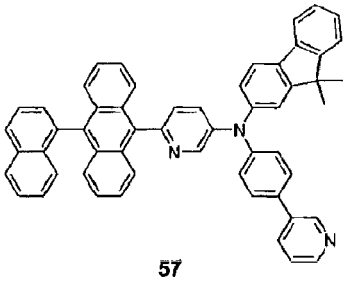

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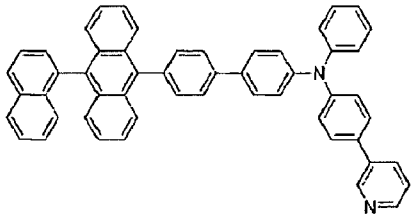

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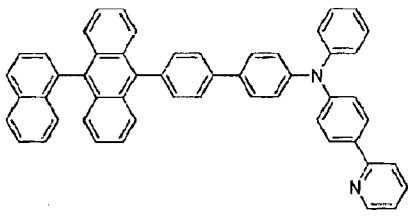

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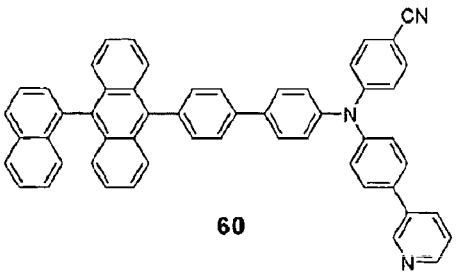

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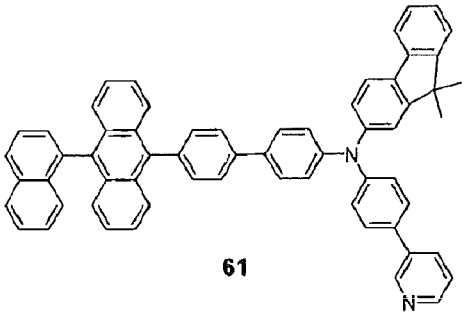

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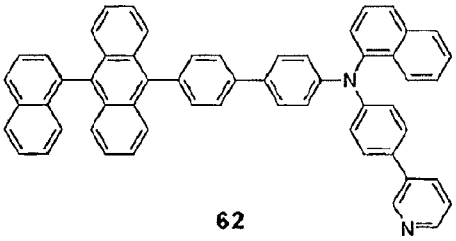

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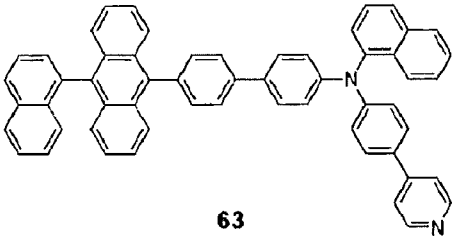

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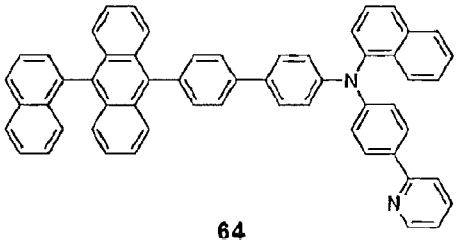

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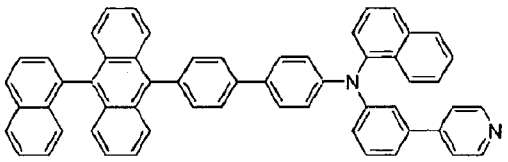

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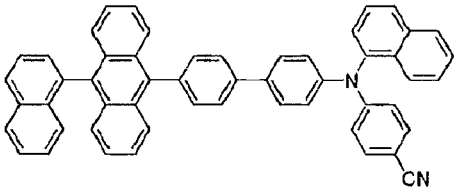

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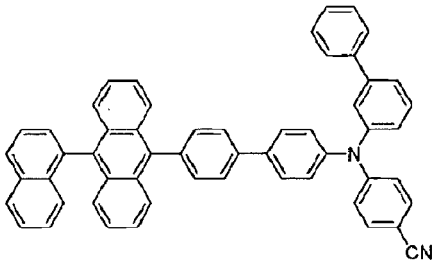

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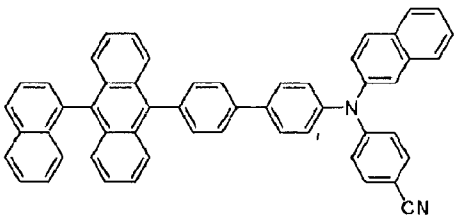

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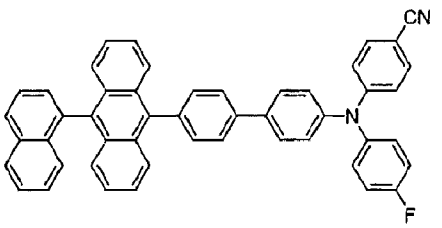

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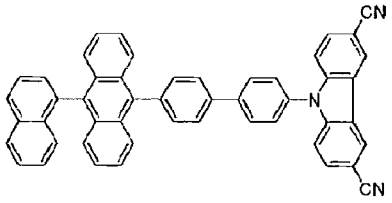

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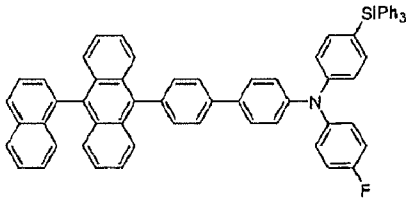

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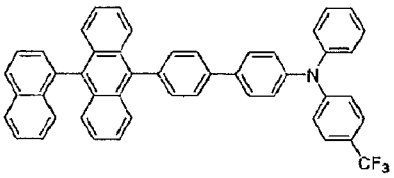

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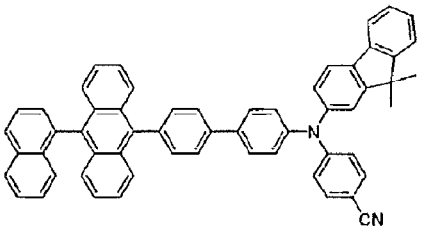

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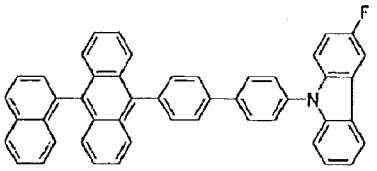

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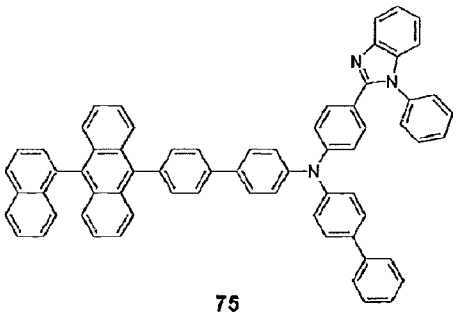

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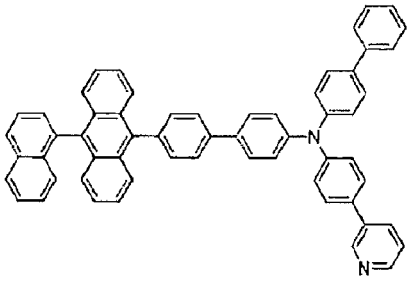

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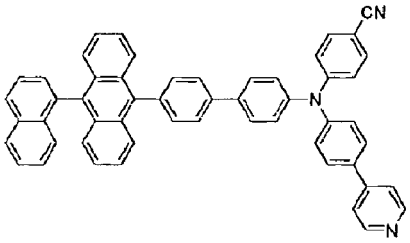

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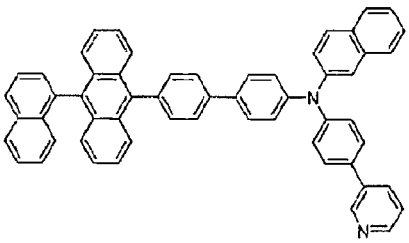

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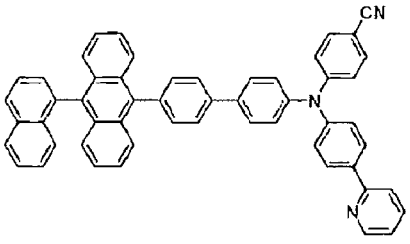

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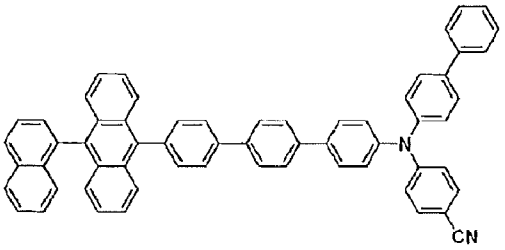

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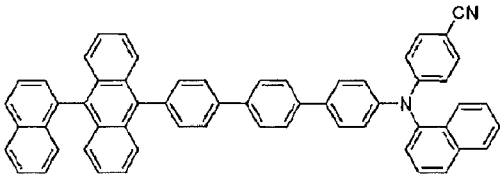

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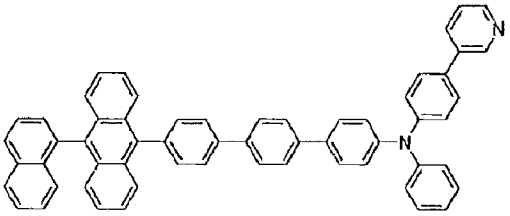

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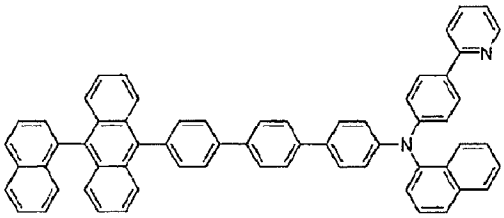

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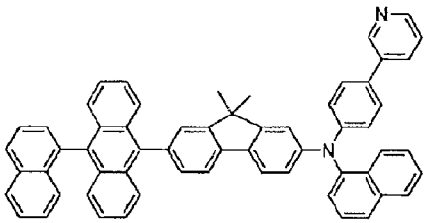

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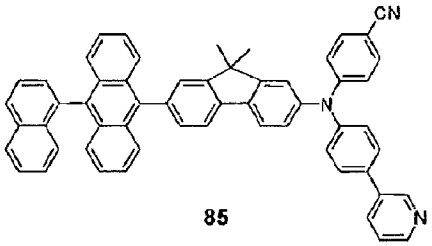

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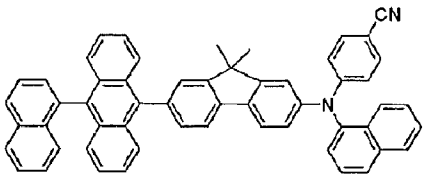

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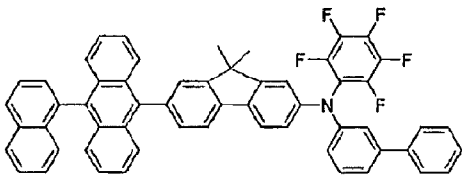

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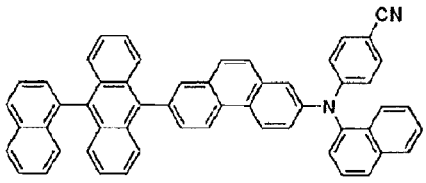

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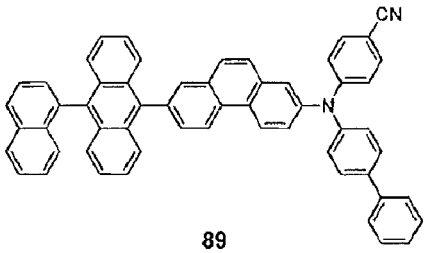

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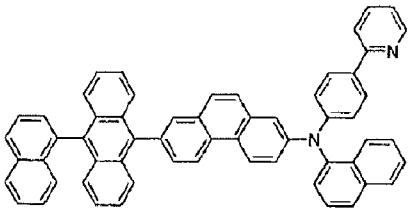

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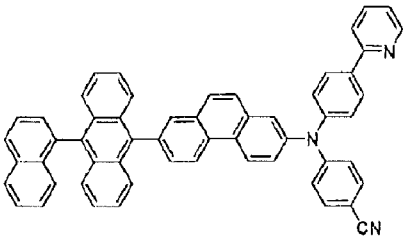

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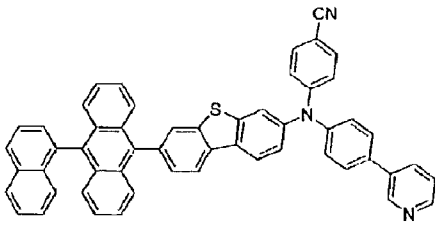

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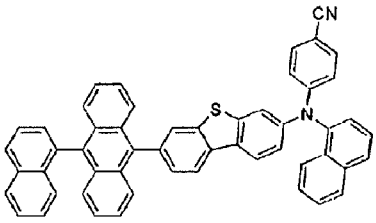

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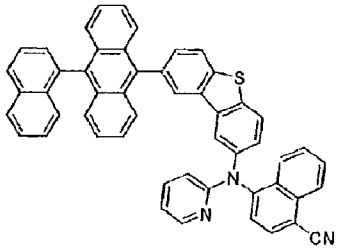

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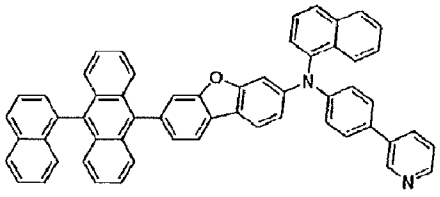

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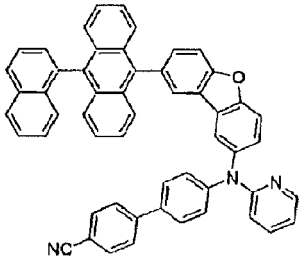

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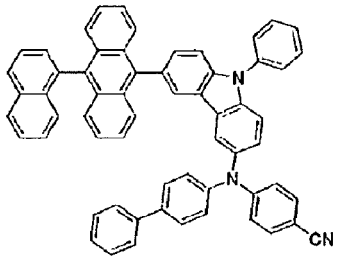

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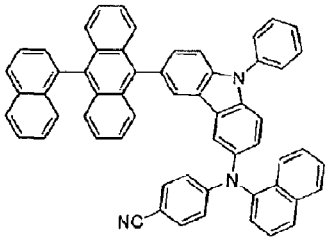

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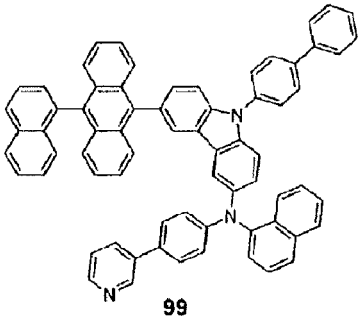

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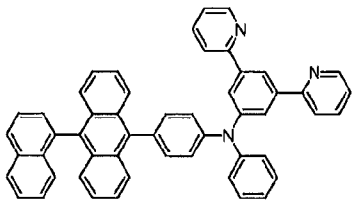

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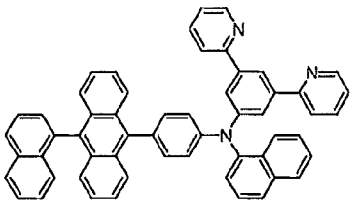

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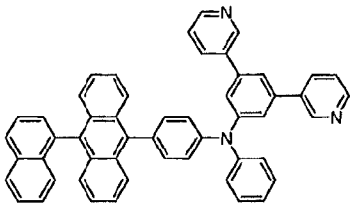

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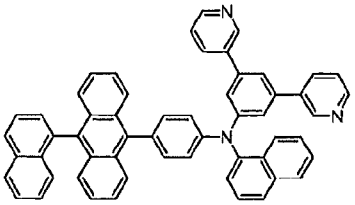

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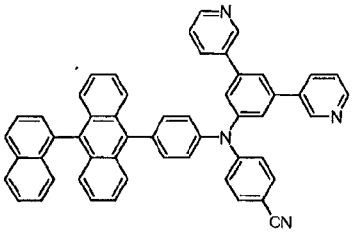

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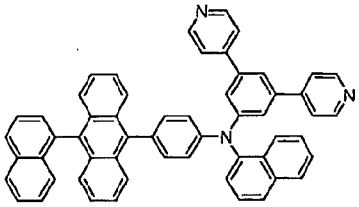

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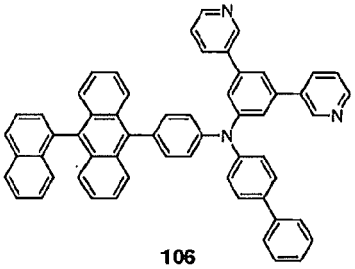

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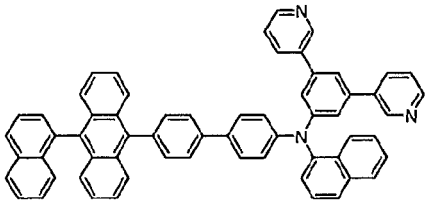

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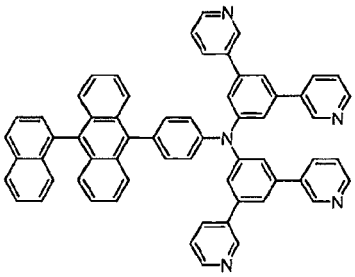

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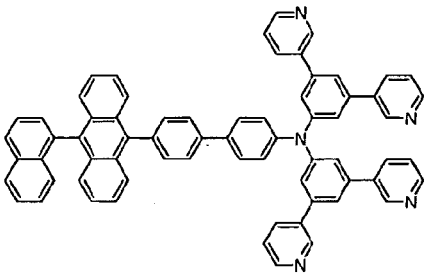

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19. Diode électroluminescente organique comprenant une première électrode, une seconde électrode placée à l'opposé de la première, et une couche organique placée entre la première et la seconde électrode, la couche organique comprenant au moins un des composés à base d'amine selon les revendications 1 à 18.

20. Diode électroluminescente organique selon la revendication 19, dans laquelle la couche organique comprend au moins l'une parmi une couche d'injection de trous, une couche de transport de trous, une couche fonctionnelle ayant à la fois des possibilités d'injection de trous et de transport de trous, une couche tampon, une couche de blocage d'électrons, une couche d'émission, une couche de blocage de trous, une couche de transport d'électrons, une couche d'injection d'électrons, et une couche fonctionnelle ayant à la fois des possibilités d'injection d'électrons et de transport d'électrons.

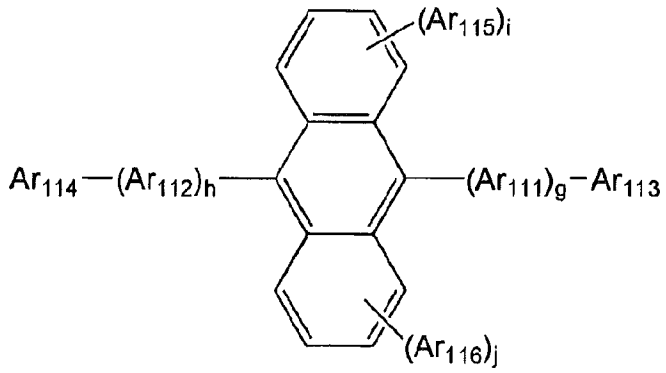
21. Diode électroluminescente organique selon la revendication 20, dans laquelle la couche organique comprend une couche de transport d'électrons, et le composé à base d'amine est compris dans la couche de transport d'électrons, de préférence,
dans laquelle la couche de transport d'électrons comprend en outre un complexe métallique.

22. Diode électroluminescente organique selon la revendication 20, dans laquelle la couche organique comprend une

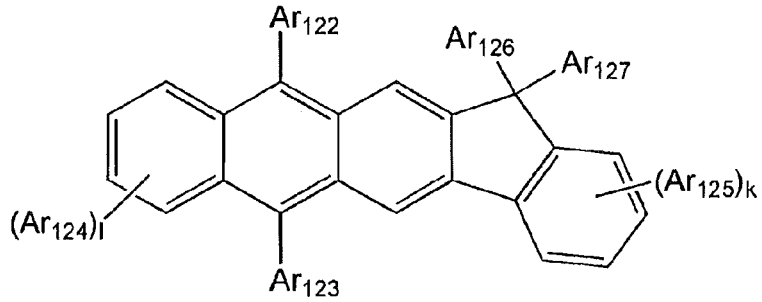
couche d'émission, et le composé à base d'amine est compris dans la couche d'émission, de préférence, dans laquelle le composé à base d'amine dans la couche d'émission joue le rôle d'hôte, et la couche d'émission comprend en outre un dopant fluorescent bleu.

23. Diode électroluminescente organique selon la revendication 22, dans laquelle le composé à base d'amine dans la couche d'émission joue le rôle de dopant, et la couche d'émission comprend en outre au moins l'un parmi un composé à base d'anthracène représenté par la Formule 400 ci-dessous et un composé à base d'anthracène représenté par la Formule 401 ci-dessous :

Formule 400



Formule 401



dans laquelle, dans les Formules 400 et 401, Ar_{111} et Ar_{112} sont chacun indépendamment un groupe arylène en C_6 à C_{60} substitué ou non ; Ar_{113} à Ar_{116} et Ar_{122} à Ar_{125} sont chacun indépendamment un groupe alkyle en C_1 à C_{10} substitué ou non, ou un groupe aryle en C_6 à C_{60} substitué ou non ; Ar_{126} et Ar_{127} sont chacun indépendamment un groupe alkyle en C_1 à C_{10} ; et g, h, i, j, k, et 1 sont chacun indépendamment un nombre entier de 0 à 4.

24. Diode électroluminescente organique selon la revendication 20, dans laquelle la couche organique comprend au moins l'une parmi une couche d'injection de trous, une couche de transport de trous, et une couche fonctionnelle ayant à la fois des possibilités d'injection de trous et de transport de trous, et l'une au moins parmi la couche d'injection de trous, la couche de transport de trous, et la couche fonctionnelle ayant à la fois des possibilités d'injection de trous et de transport de trous comprend un dopant p.

REFERENCES CITED IN THE DESCRIPTION

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Non-patent literature cited in the description

- Synthesis and electroluminescent properties of blue fluorescent triphenylamine substituted anthracene derivatives for OLEDs. *Molecular Crystals and Liquid Crystals,* 11 October 2010, vol. 530 (1 **[0006]**

专利名称(译)	胺类化合物和包括其的有机发光二极管		
公开(公告)号	EP2626399B1	公开(公告)日	2014-12-03
申请号	EP2013154326	申请日	2013-02-07
[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.		
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IPC分类号	C09K11/06 C07C15/28 H01L51/00 H05B33/10		
CPC分类号	C09K11/06 C07C211/56 C07C211/59 C07C211/61 C07C217/92 C07C255/58 C07C2603/18 C07C2603/24 C07C2603/26 C07C2603/50 C07D209/88 C07D213/36 C07D213/74 C07D307/91 C07D333/76 C07F7/0805 C09K2211/1007 C09K2211/1011 C09K2211/1014 H01L51/0058 H01L51/006 H01L51/0081 H05B33/10		
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优先权	1020120012532 2012-02-07 KR 1020120061676 2012-06-08 KR		
其他公开文献	EP2626399A1		
外部链接	Espacenet		

摘要(译)

胺类化合物和包含胺类化合物的有机发光二极管。

Formula 1