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(19) **United States**(12) **Patent Application Publication**
SHIN et al.(10) **Pub. No.: US 2014/0291646 A1**(43) **Pub. Date: Oct. 2, 2014**(54) **PYRENE-BASED COMPOUND AND
ORGANIC LIGHT-EMITTING DIODE
INCLUDING THE SAME**(52) **U.S. CL.**
CPC *H01L 51/006* (2013.01); *H01L 51/0061*
(2013.01)USPC **257/40**; 546/304; 544/330; 544/292;
544/326; 544/356(71) Applicant: **Samsung Display Co., Ltd.**,
Yongin-City (KR)(72) Inventors: **Dong-Woo SHIN**, Yongin-City (KR);
O-Hyun KWON, Yongin-City (KR);
Mie-Hwa PARK, Yongin-City (KR);
Seul-Ong KIM, Yongin-City (KR);
Byoung-Ki CHOI, Yongin-City (KR)(57) **ABSTRACT**

A pyrene-based compound and an organic light-emitting diode including the same, the pyrene-based compound being represented by Formula 1, below:

(73) Assignee: **Samsung Display Co., Ltd.**,
Yongin-City (KR)(21) Appl. No.: **14/224,679**(22) Filed: **Mar. 25, 2014**(30) **Foreign Application Priority Data**

Mar. 27, 2013 (KR) 10-2013-0033082

Publication Classification(51) **Int. Cl.**
H01L 51/00 (2006.01)

<Formula 1>

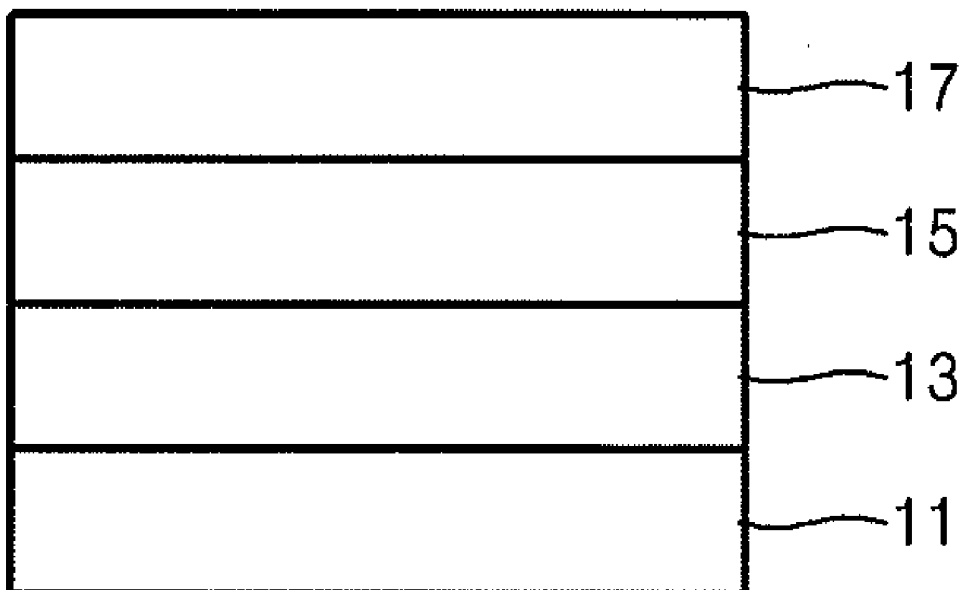
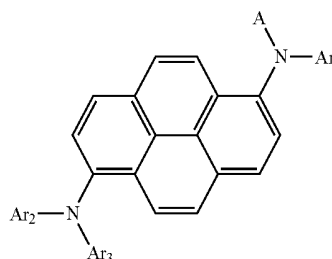
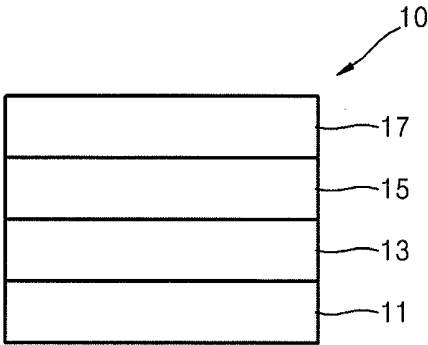


FIG. 1



**PYRENE-BASED COMPOUND AND
ORGANIC LIGHT-EMITTING DIODE
INCLUDING THE SAME**

**CROSS-REFERENCE TO RELATED
APPLICATION**

[0001] Korean Patent Application No. 10-2013-0033082, filed on Mar. 27, 2013, in the Korean Intellectual Property Office, and entitled: "Pyrene-Based Compound and Organic Light-Emitting Diode Including the Same," is incorporated by reference herein in its entirety.

BACKGROUND

[0002] 1. Field

[0003] Embodiments relates to a pyrene-based compound and an organic light-emitting diode including the same.

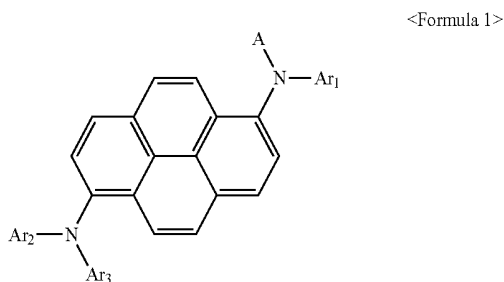
[0004] 2. Description of the Related Art

[0005] Organic light-emitting diodes (OLEDs), which are self-emitting devices, may have advantages such as wide viewing angles, excellent contrast, quick response, high brightness, excellent driving voltage characteristics, and a may provide multicolored images.

SUMMARY

[0006] Embodiments are directed to a pyrene-based compound and an organic light-emitting diode including the same.

[0007] The embodiments may be realized by providing a pyrene-based compound represented by Formula 1, below:

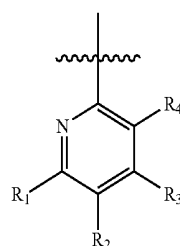


[0008] wherein, in Formula 1, A is a substituted or unsubstituted C₂-C₃₀ heteroaromatic group including at least one nitrogen (N); and Ar₁, Ar₂, and Ar₃ are each independently a substituted or unsubstituted phenyl group or a substituted or unsubstituted naphthyl group.

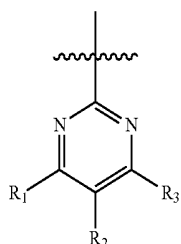
[0009] A may be selected from a substituted or unsubstituted pyridinyl group, a substituted or unsubstituted pyrimidinyl group, a substituted or unsubstituted triazinyl group, a substituted or unsubstituted pyrazinyl group, a substituted or unsubstituted pyridazinyl group, a substituted or unsubstituted isoindolyl group, a substituted or unsubstituted indolyl group, a substituted or unsubstituted indazolyl group, a substituted or unsubstituted isoquinolinyl group, a substituted or unsubstituted quinolinyl group, a substituted or unsubstituted phthalazinyl group, a substituted or unsubstituted quinoxalinyl group, a substituted or unsubstituted quinazolinyl group, a substituted or unsubstituted cinnolinyl group, and a substituted or unsubstituted 1,10-phenanthrolinyl group.

[0010] A may be selected from a pyridinyl group, a pyrimidinyl group, a triazinyl group, a pyrazinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, an isoquinolinyl group, a quinolinyl group, a phthalazinyl group, a quinoxalinyl group, a cinnolinyl group, and a 1,10-phenanthrolinyl group; or a pyridinyl group, a pyrimidinyl group, a triazinyl group, a pyrazinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, an isoquinolinyl group, a quinolinyl group, a phthalazinyl group, a quinoxalinyl group, a cinnolinyl group, and a 1,10-phenanthrolinyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof, a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, and a C₂-C₆₀ heteroaryl group, a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, and a C₂-C₆₀ heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group, or —Si(Q₁₁)(Q₁₂)(Q₁₃), in which Q₁₁ to Q₁₃ are each independently a C₁-C₆₀ alkyl group, a C₁-C₆₀ alkoxy group, a C₆-C₆₀ aryl group, or a C₂-C₆₀ heteroaryl group.

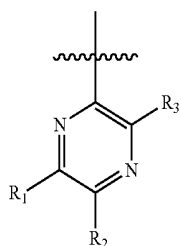
[0011] A may be a group represented by one of Formulae 2A to 2K below:



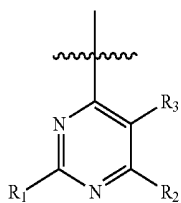
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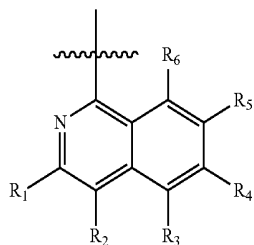
Formula 2B



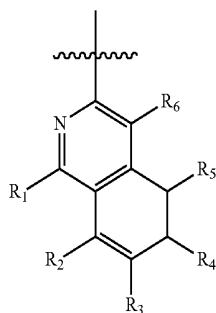
Formula 2C



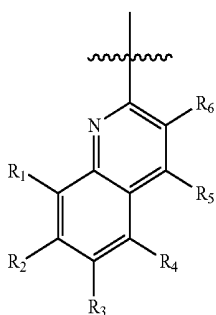
Formula 2D



Formula 2E

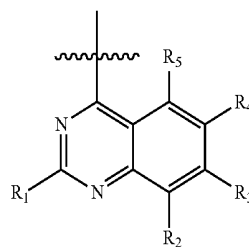


Formula 2F

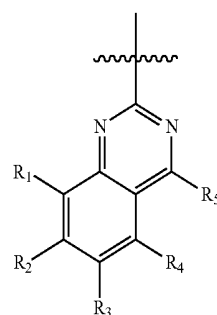


Formula 2G

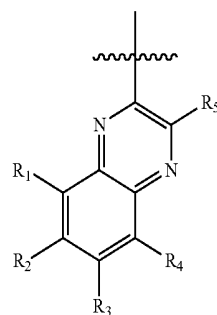
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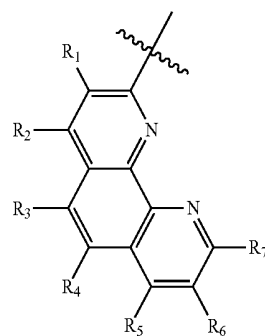
Formula 2H



Formula 2I



Formula 2J



Formula 2K

[0012] wherein, in Formulae 2A to 2K, R_1 to R_7 are each independently selected from a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group; a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof; a C_3 - C_{10} cycloalkyl group, a C_3 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_3 - C_{10} heterocycloalkenyl group, a

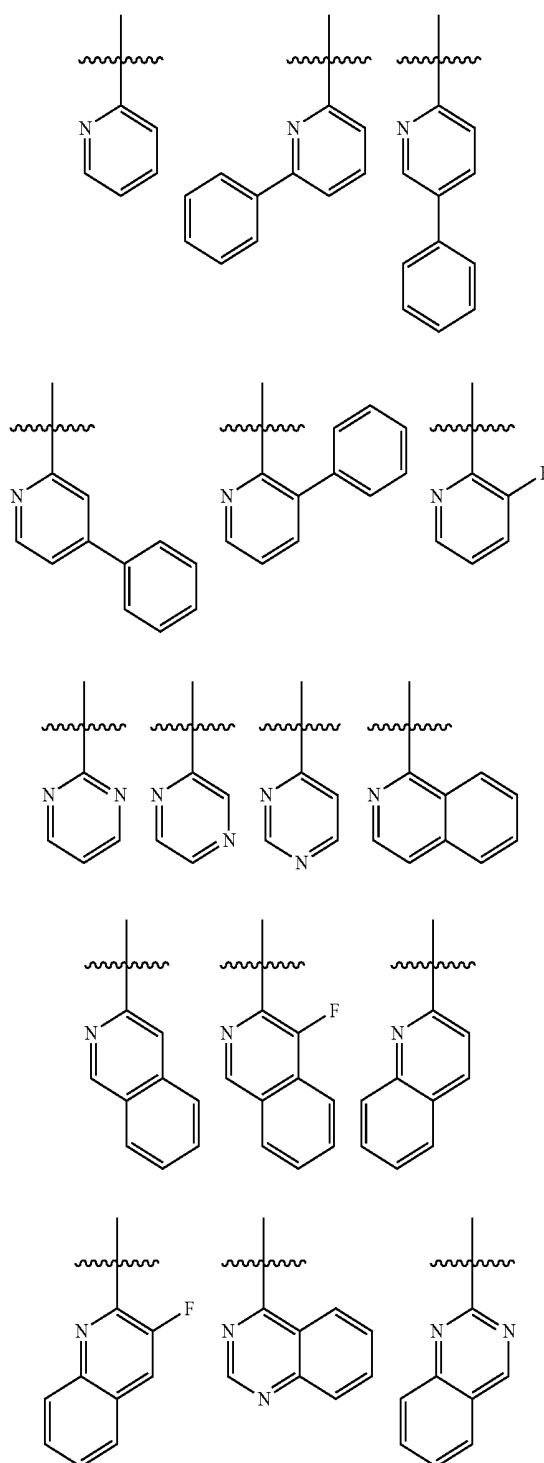
C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, and a C₂-C₆₀ heteroaryl group; a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, and a C₂-C₆₀ heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group; and —Si(Q₁₁)(Q₁₂)(Q₁₃), in which Q₁₁ to Q₁₃ are each independently a C₁-C₆₀ alkyl group, a C₁-C₆₀ alkoxy group, a C₆-C₆₀ aryl group, or a C₂-C₆₀ heteroaryl group.

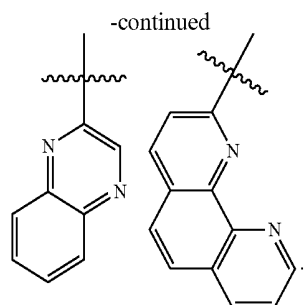
[0013] R₁ to R₇ may each independently be selected from a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a C₁-C₁₀ alkyl group; a C₁-C₆₀ alkyl group substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof; a C₆-C₁₆ aryl group, and a C₂-C₁₆ heteroaryl group; a C₆-C₁₆ aryl group or a C₂-C₁₆ heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₁₀ alkyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group; and —Si(Q₁₁)(Q₁₂)(Q₁₃), in which Q₁₁ to Q₁₃ are each independently a C₁-C₁₀ alkyl group, a C₆-C₁₆ aryl group, or a C₂-C₁₆ heteroaryl group.

[0014] R₁ to R₇ may each independently be selected from a hydrogen atom, a deuterium atom, F, Cl, Br, I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid 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, and a hexyl group; a phenyl group, a naphthyl group, and an anthracenyl group; and a phenyl group, a naphthyl group, and an anthracenyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a methyl group,

an ethyl group, a propyl group, a butyl group, a pentyl group, a hexyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, and a phenylcarbazolyl group.

[0015] A may have a structure selected from one of the following structures:



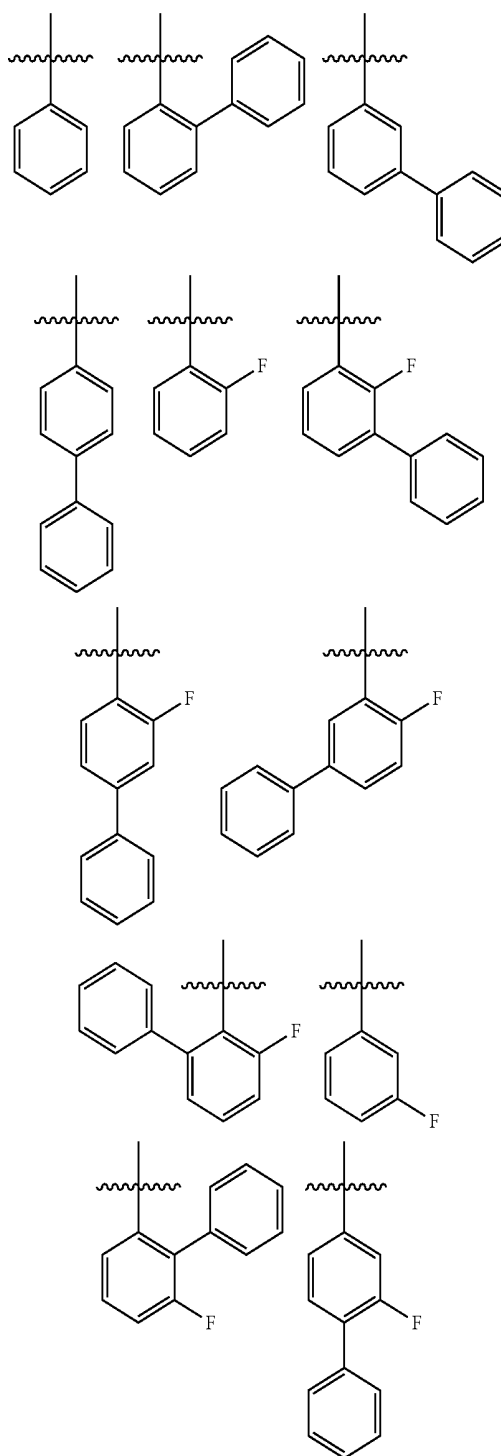


[0016] Ar_1 , Ar_2 , and Ar_3 may each independently be selected from a phenyl group and a naphthyl group; and a phenyl group and a naphthyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group, a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof, a C_3 - C_{10} cycloalkyl group, a C_3 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_3 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, and a C_2 - C_{60} heteroaryl group, a C_3 - C_{10} cycloalkyl group, a C_3 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_3 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, and a C_2 - C_{60} heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic 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 anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group, and $-Si(Q_{11})(Q_{12})(Q_{13})$, in which Q_{11} to Q_{13} are each independently a C_1 - C_{60} alkyl group, a C_1 - C_{60} alkoxy group, a C_6 - C_{60} aryl group, or a C_2 - C_{60} heteroaryl group.

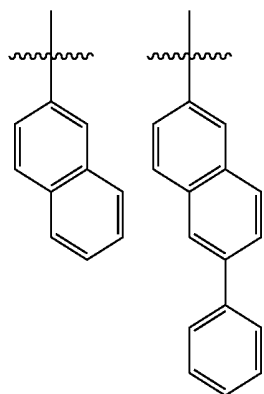
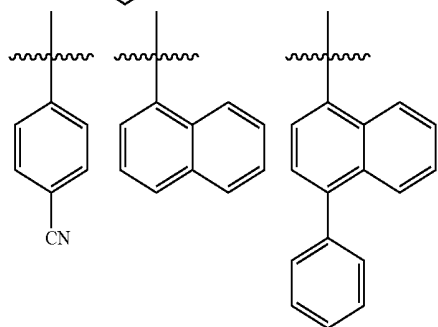
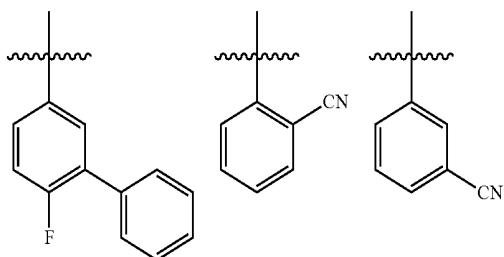
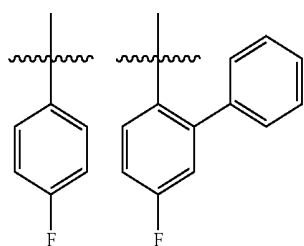
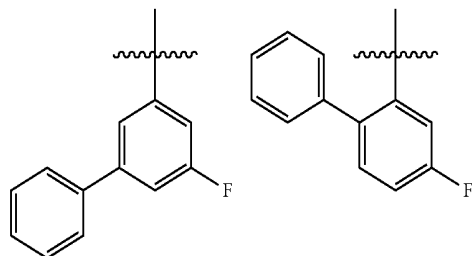
[0017] Ar_1 , Ar_2 , and Ar_3 may each independently be selected from a phenyl group and a naphthyl group; and a phenyl group and a naphthyl group, each substituted with at least one of a deuterium atom, F, Cl, Br, I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof and a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, and a hexyl group, a phenyl group, a naphthyl group, and an anthracenyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a

hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, a hexyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, and a phenylcarbazolyl group.

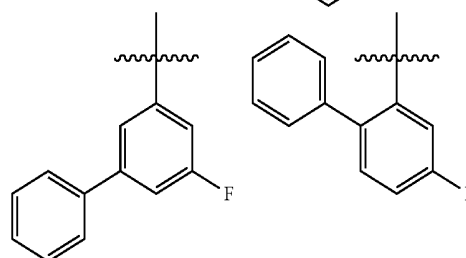
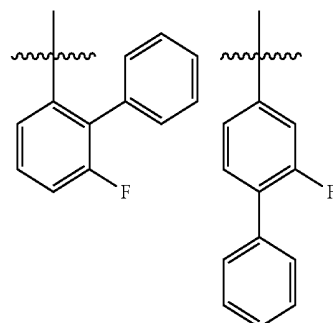
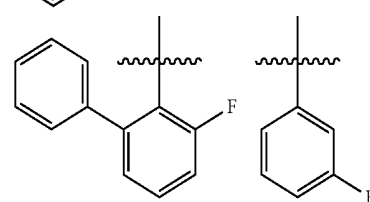
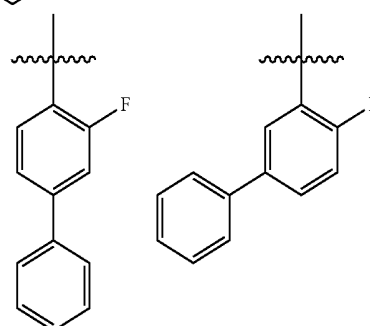
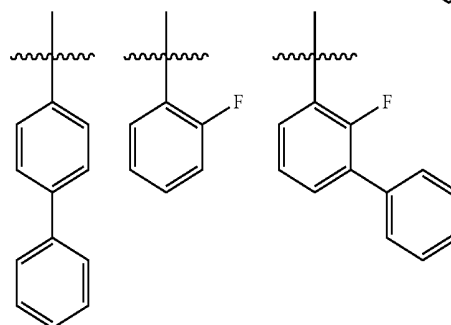
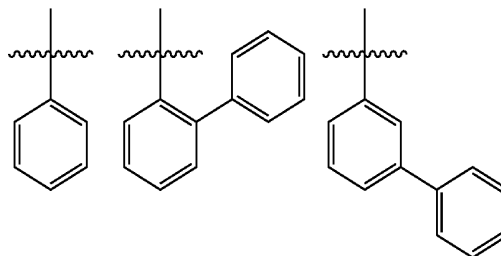
[0018] Ar_1 may be selected from one of the following structures:



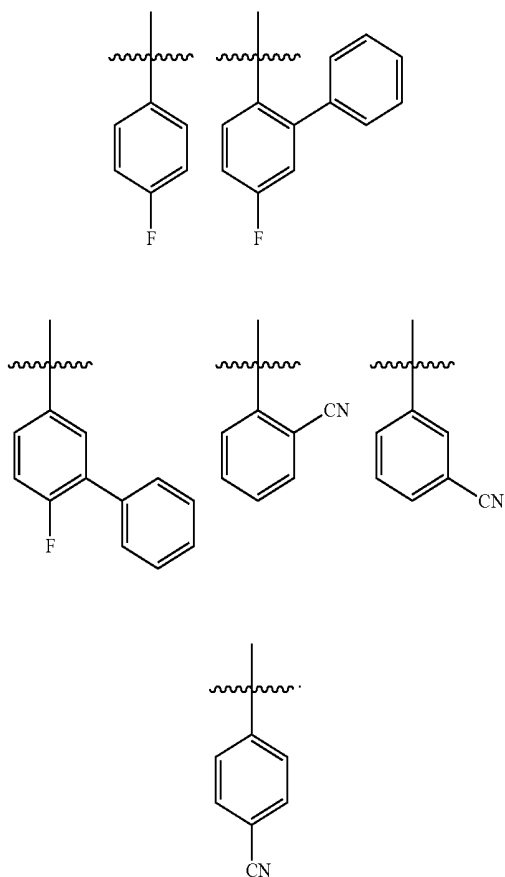
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[0019] Ar_2 and Ar_3 may each independently be selected from one of the following structures:

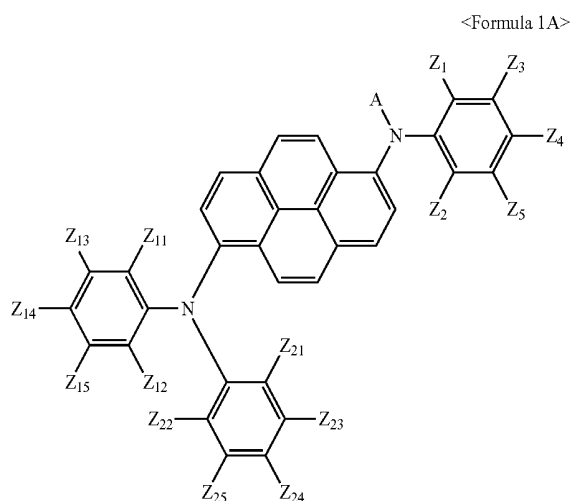


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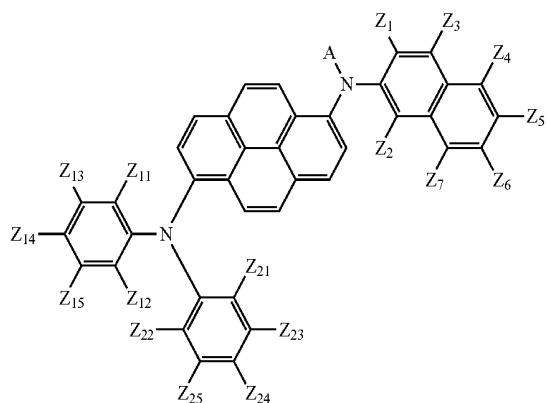
[0020] Ar₁ may be a substituted or unsubstituted phenyl group or a substituted or unsubstituted naphthyl group, and Ar₂ and Ar₃ are a substituted or unsubstituted phenyl group.

[0021] The pyrene-based compound may be represented by one of Formulae 1A, 1B, or 1C, below:

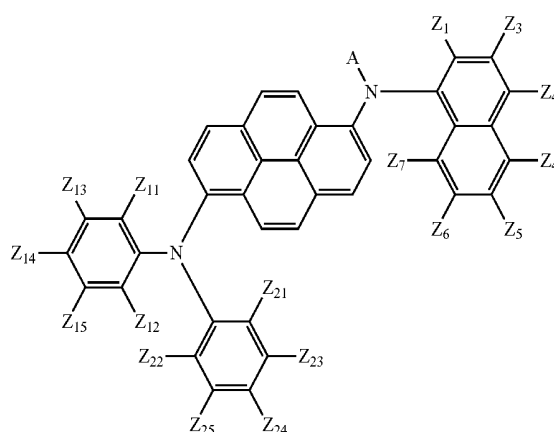


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<Formula 1B>

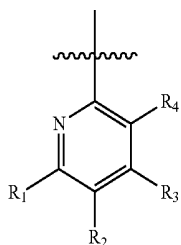


<Formula 1C>

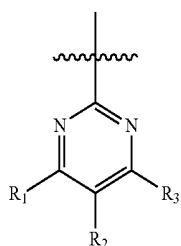


[0022] wherein, in Formulae 1A, 1B, and 1C, A is a substituted or unsubstituted C₂-C₃₀ heteroaromatic group including at least one nitrogen (N); and Z₁ to Z₇, Z₁₁ to Z₁₅ and Z₂₁ to Z₂₅ are each independently selected from a hydrogen atom, a deuterium atom, F, Cl, Br, I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid 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, and a hexyl group; a phenyl group, a naphthyl group, and an anthracenyl group; and a phenyl group, a naphthyl group, and an anthracenyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, a hexyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group and a phenylcarbazolyl group.

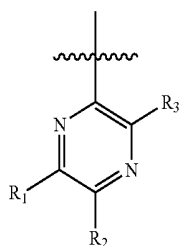
[0023] A may be represented by one of Formulae 2A to 2K, below:



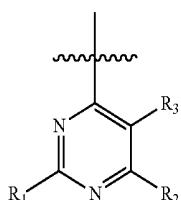
Formula 2A



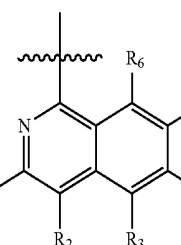
Formula 2B



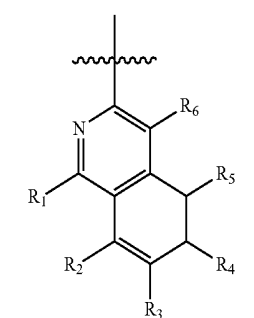
Formula 2C



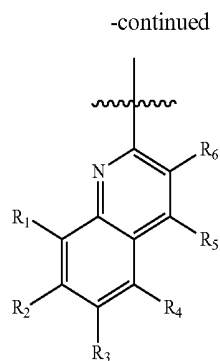
Formula 2D



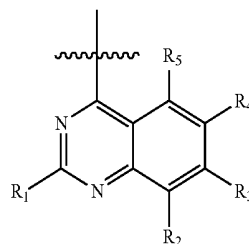
Formula 2E



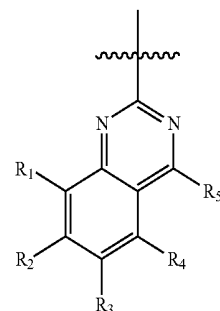
Formula 2F



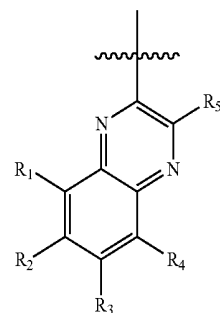
Formula 2G



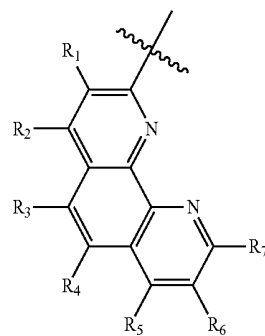
Formula 2H



Formula 2I



Formula 2J



Formula 2K

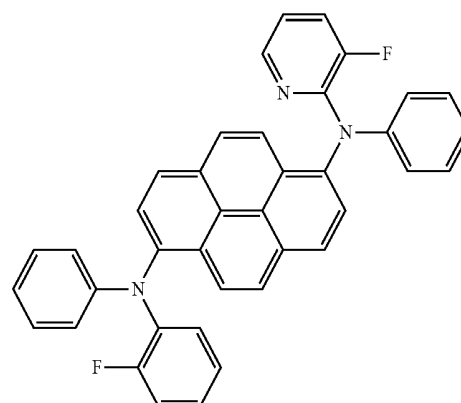
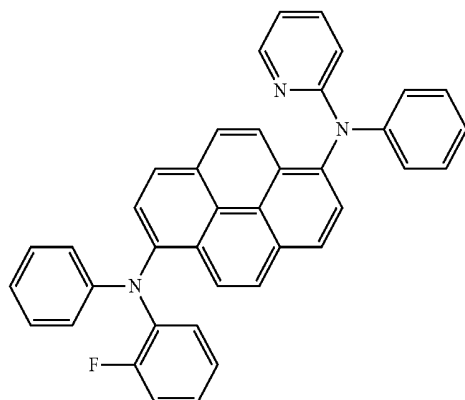
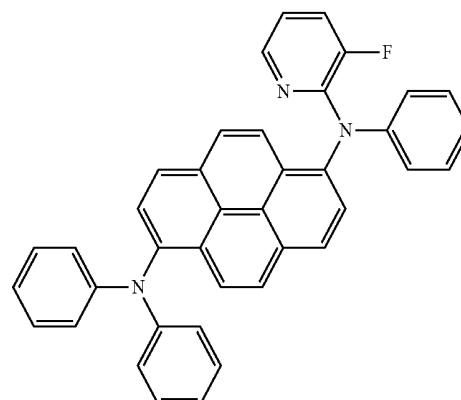
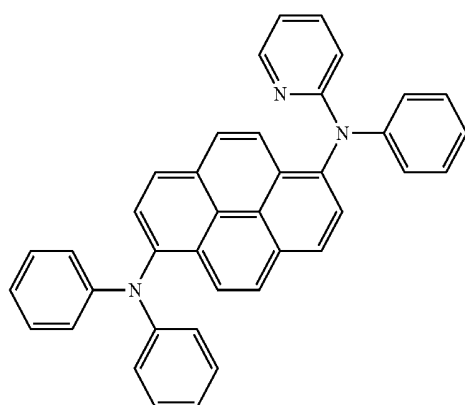
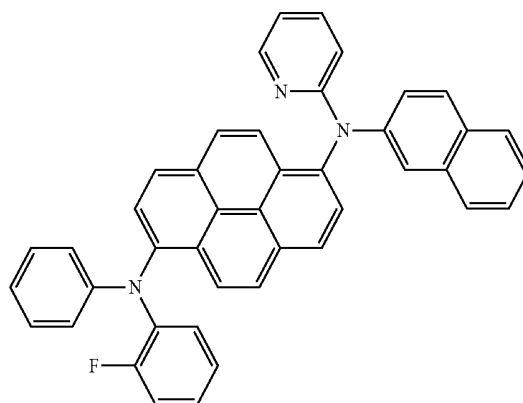
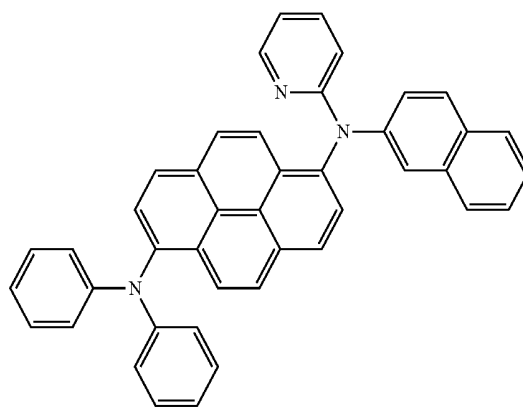
[0024] wherein, in Formulae 2A to 2K, R₁ to R₇ are each independently selected from a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine

group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a C_1 - C_{10} alkyl group; a C_1 - C_{60} alkyl group substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof; a C_6 - C_{16} aryl group and a C_2 - C_{16} heteroaryl group; a C_6 - C_{16} aryl group and a C_2 - C_{16} heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C_1 - C_{10} alkyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group; and $-\text{Si}(\text{Q}_{11})(\text{Q}_{12})(\text{Q}_{13})$, in which Q_{11} to Q_{13} are each independently a C_1 - C_{10} alkyl group, a C_6 - C_{16} aryl group, or a C_2 - C_{16} heteroaryl group.

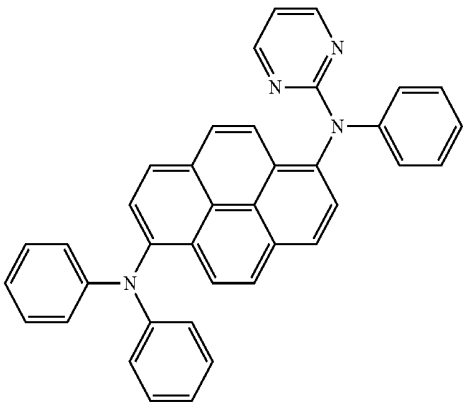
[0025] R_1 to R_7 , Z_1 to Z_7 , Z_{11} to Z_{15} and Z_{21} to Z_{25} may each independently be a hydrogen atom, a F atom, a cyano group, or a phenyl group.

[0026] The pyrene-based compound may be one of Compounds I to 96, below:

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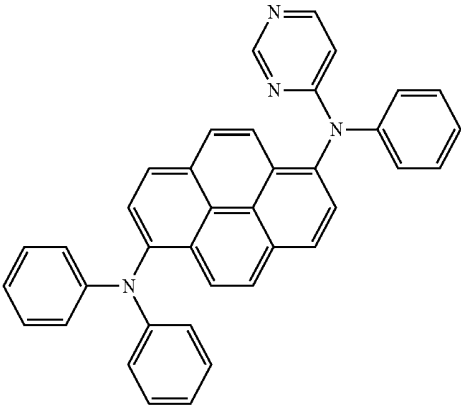


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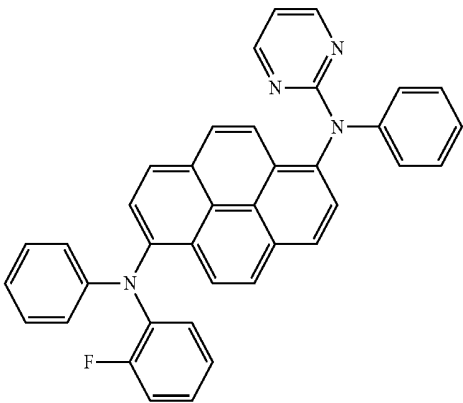


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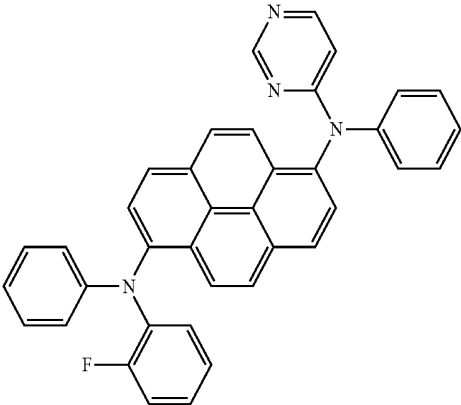
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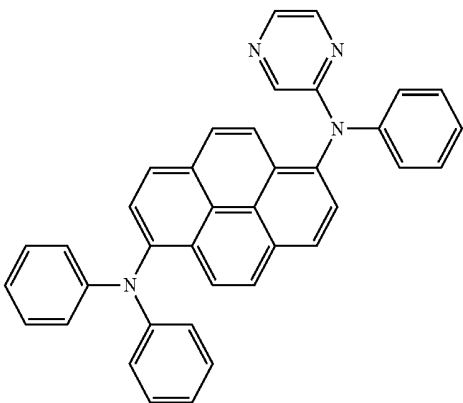
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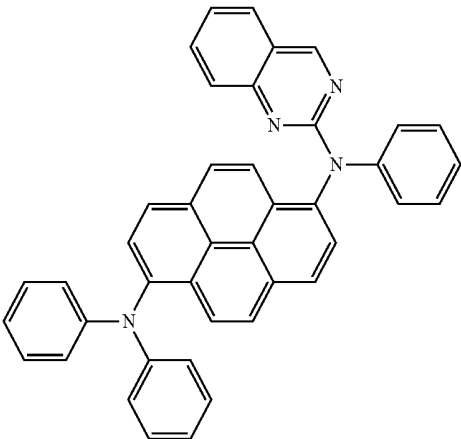
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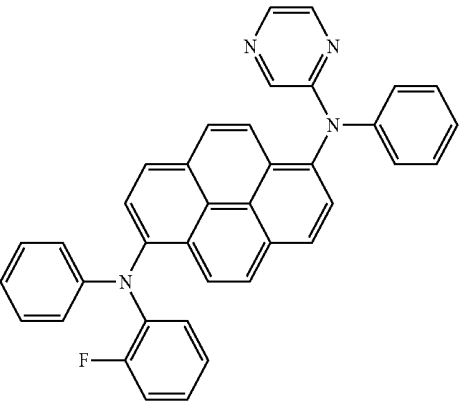
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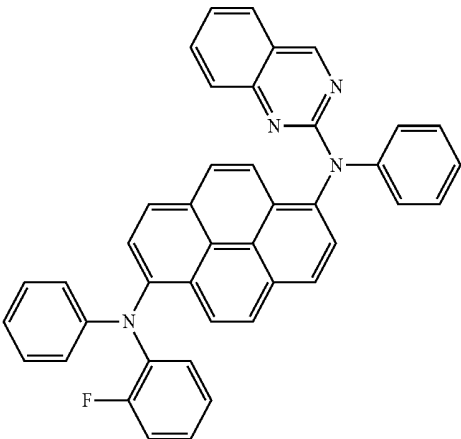
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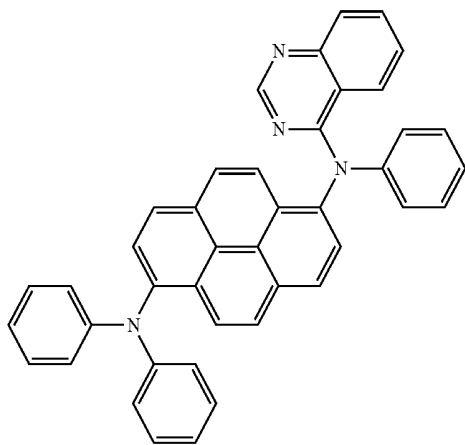


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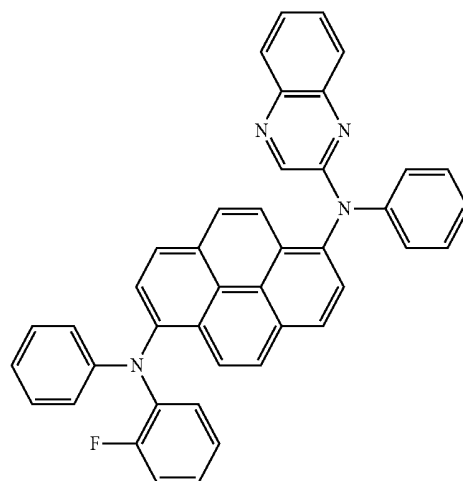
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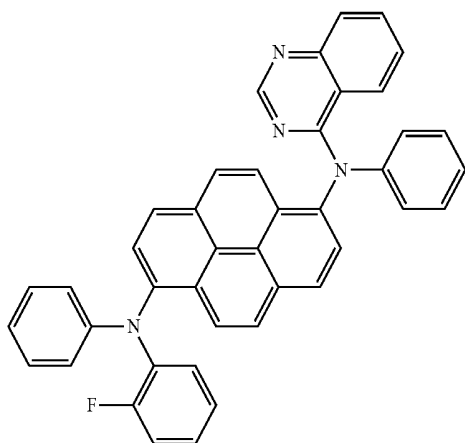


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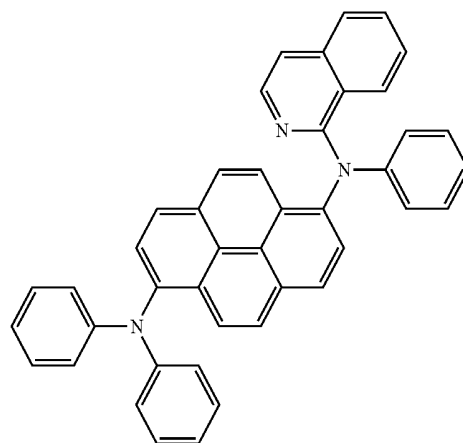
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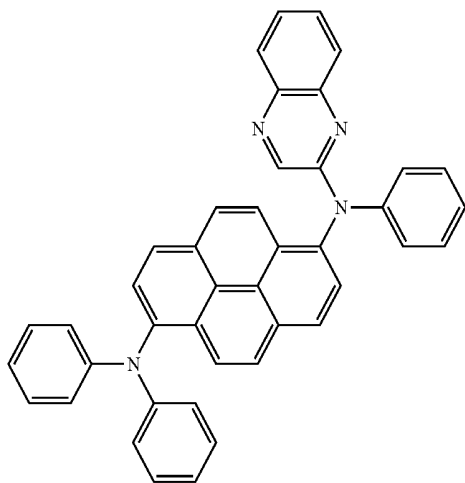
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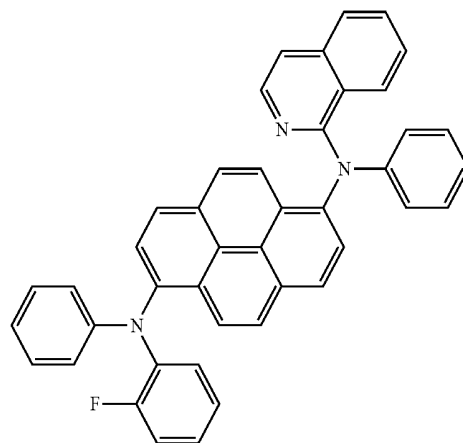
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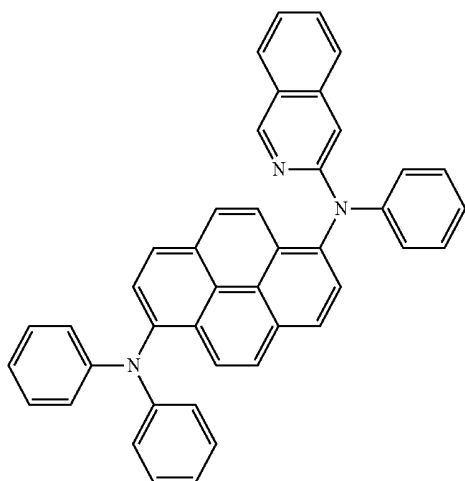


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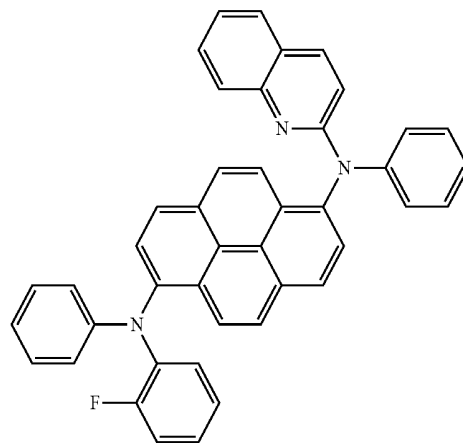
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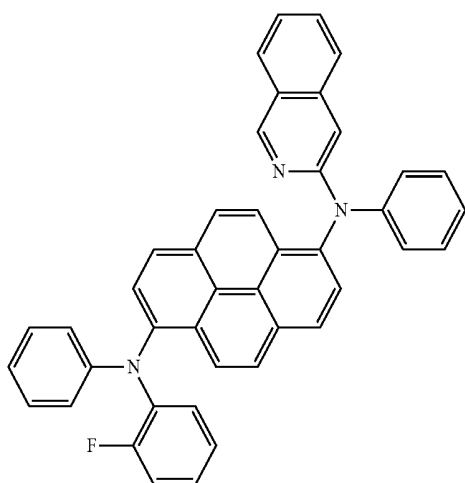


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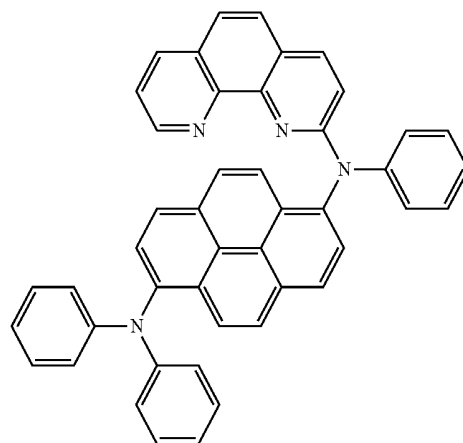
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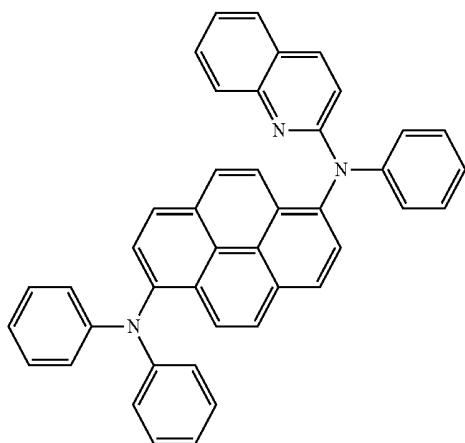
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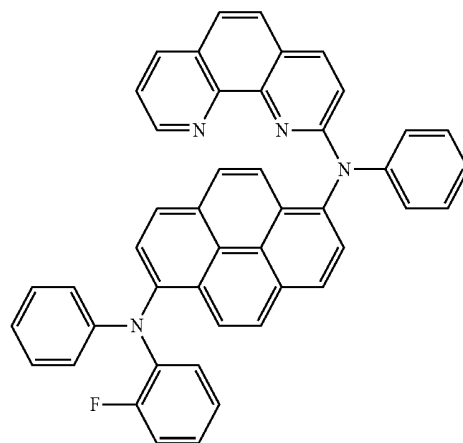
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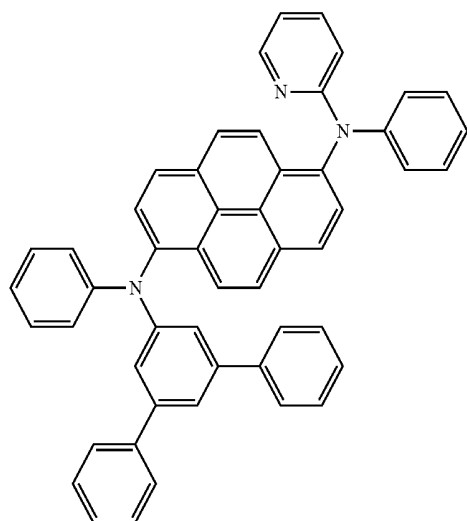


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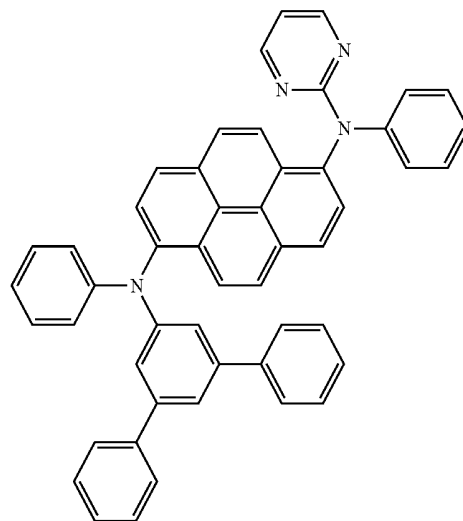
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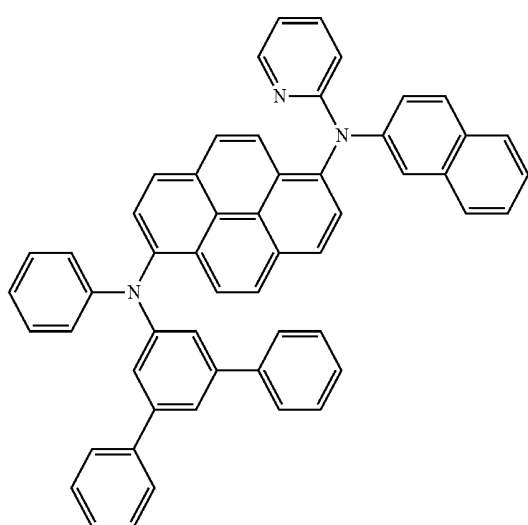


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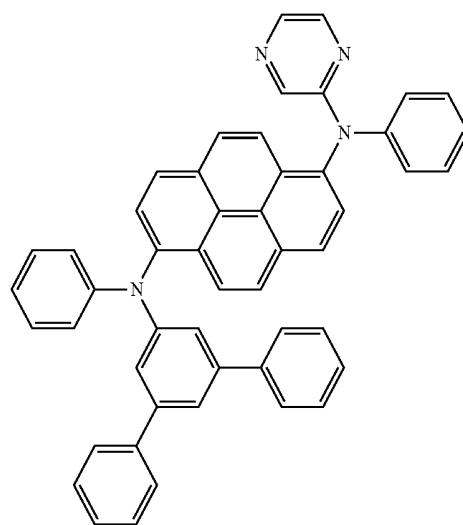
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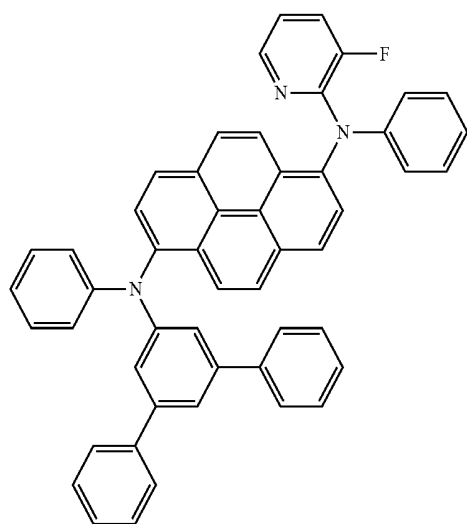
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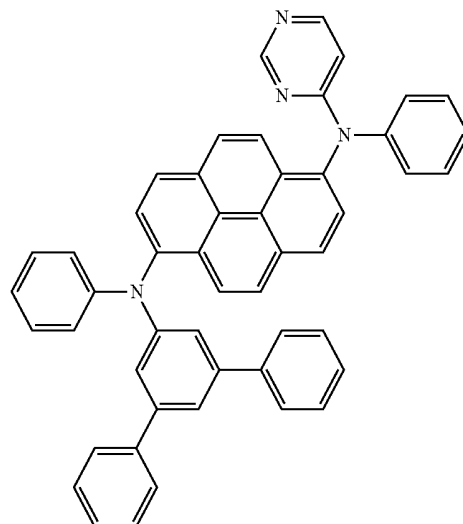
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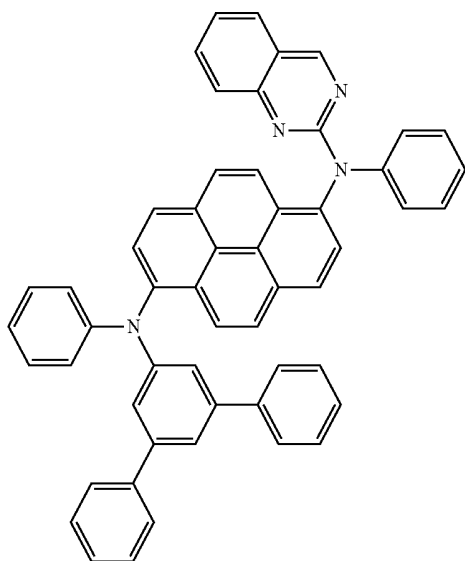
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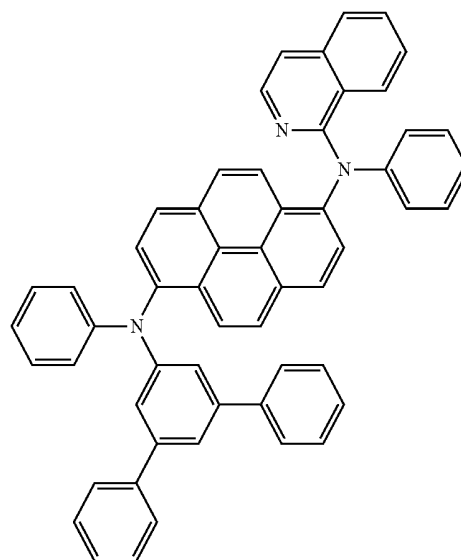
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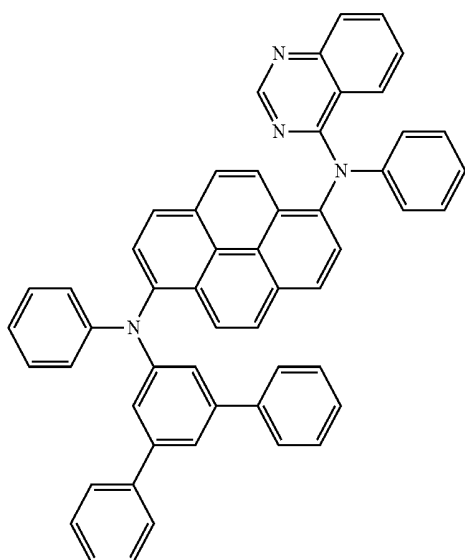


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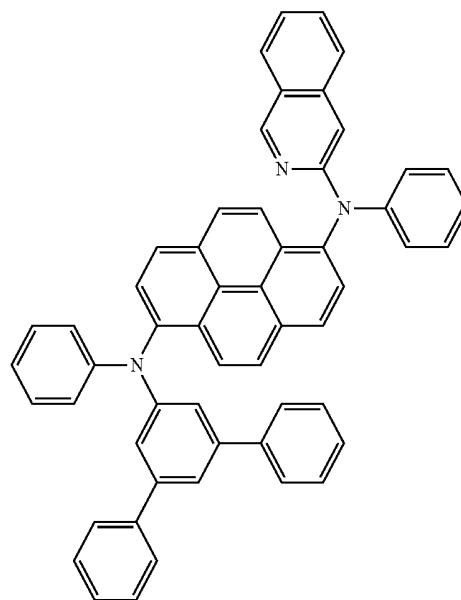
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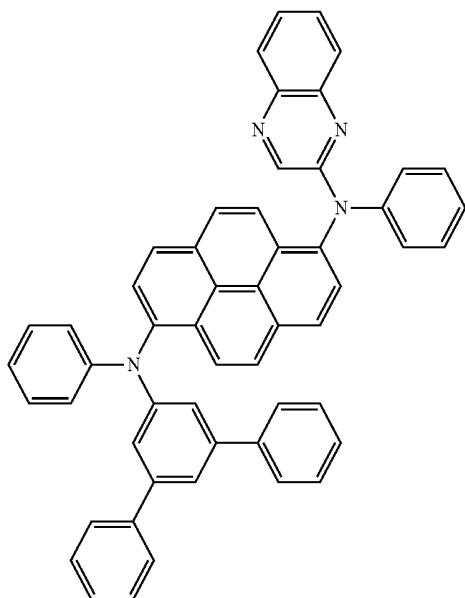
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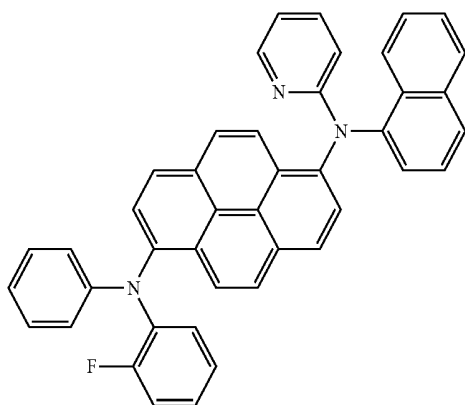
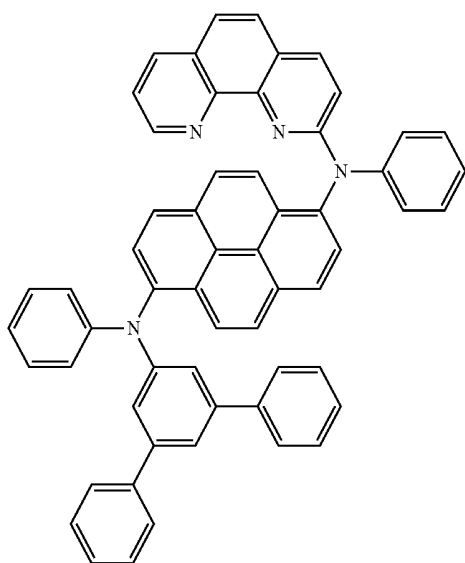
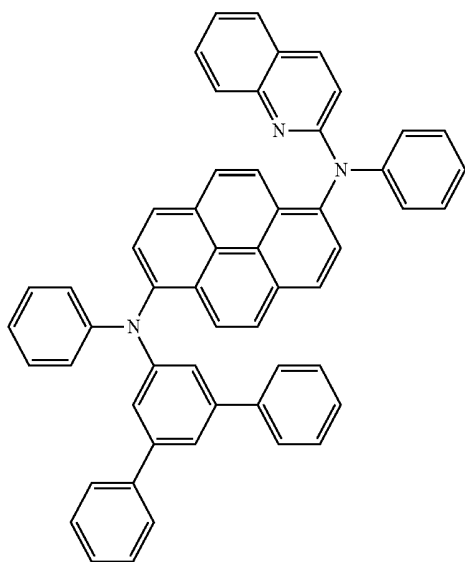
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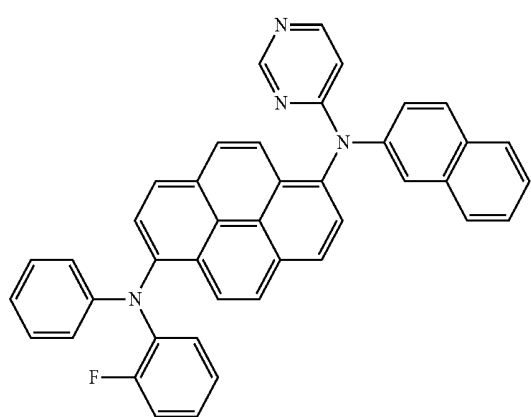
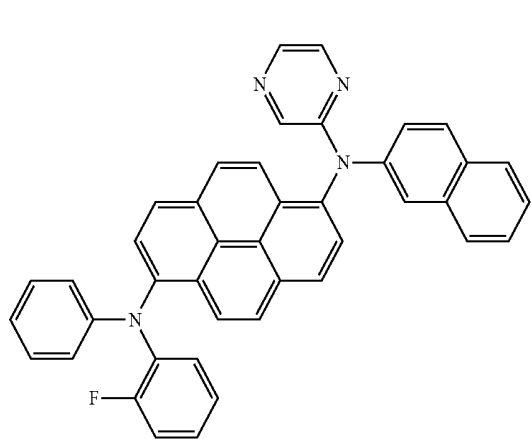
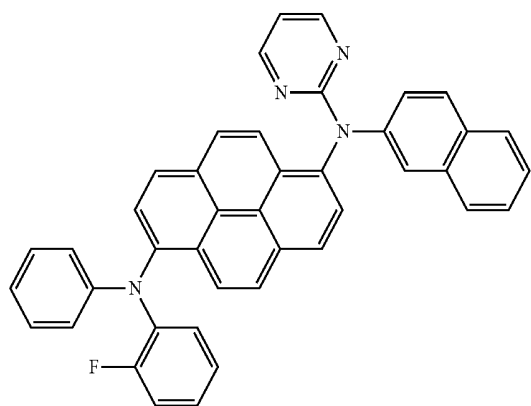
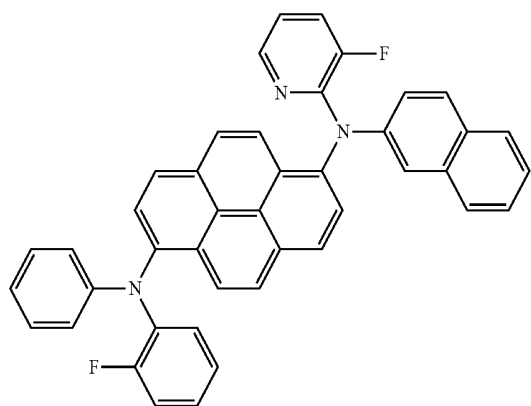
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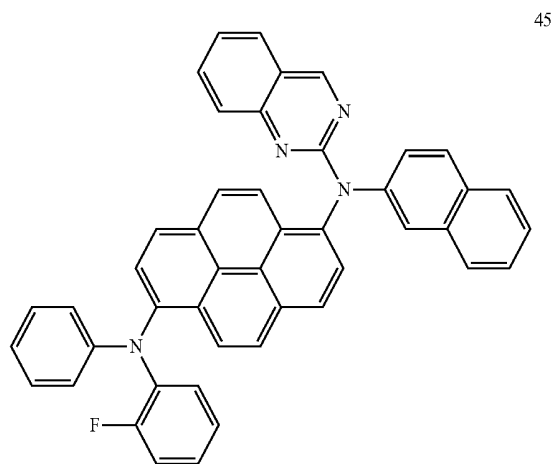
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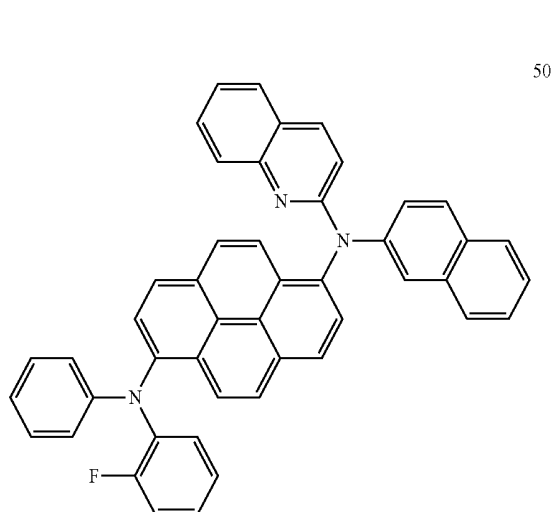
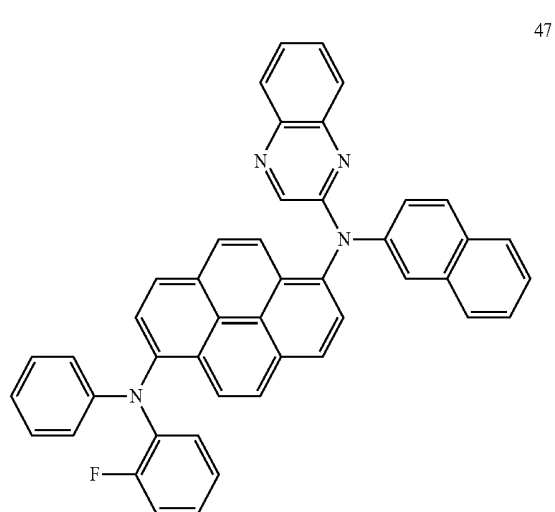
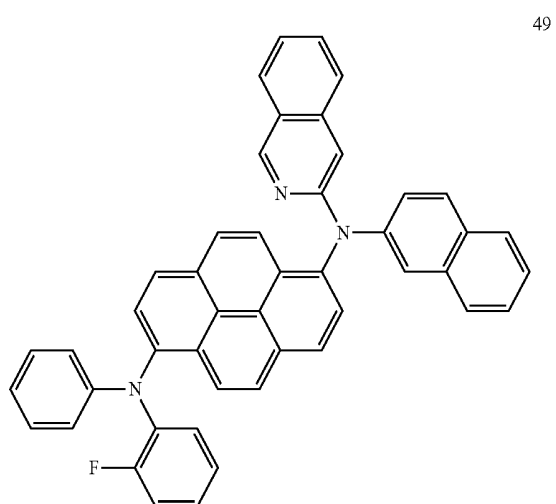
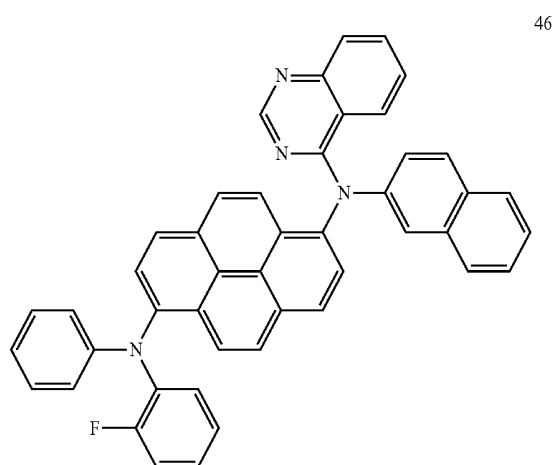
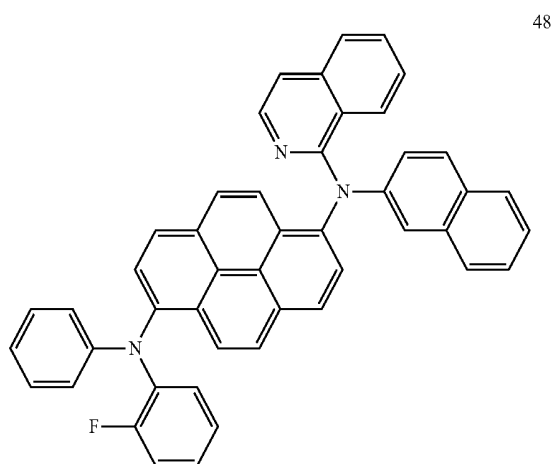
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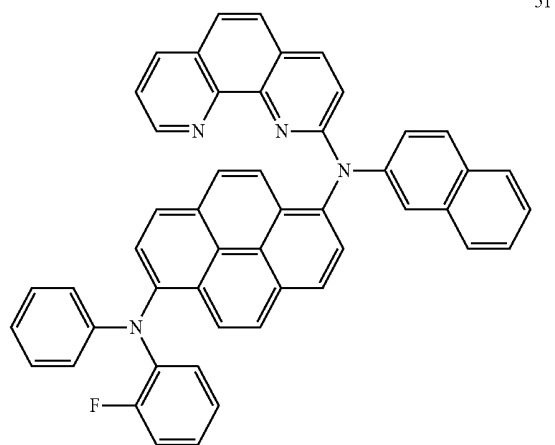
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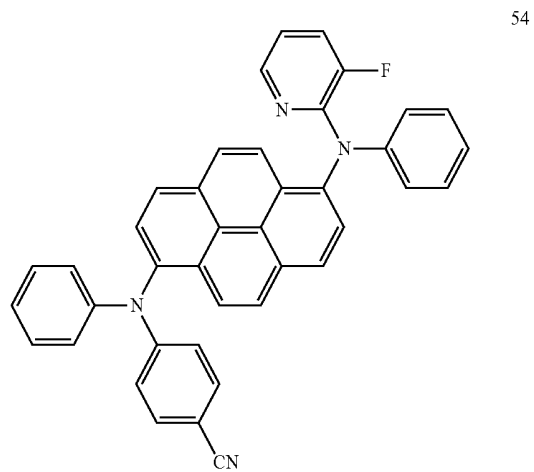
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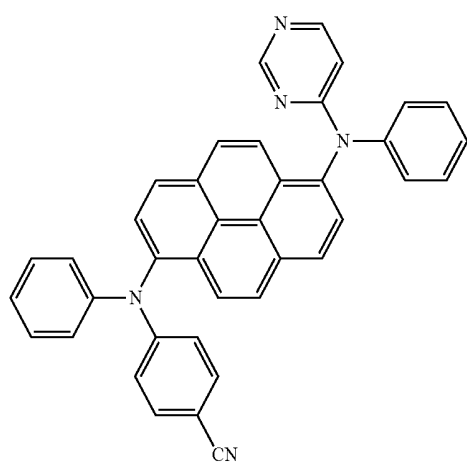
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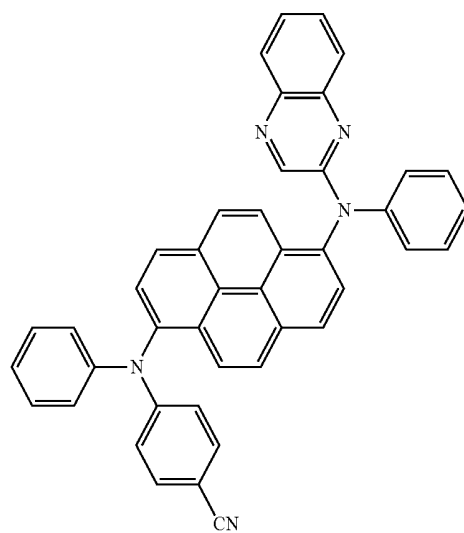
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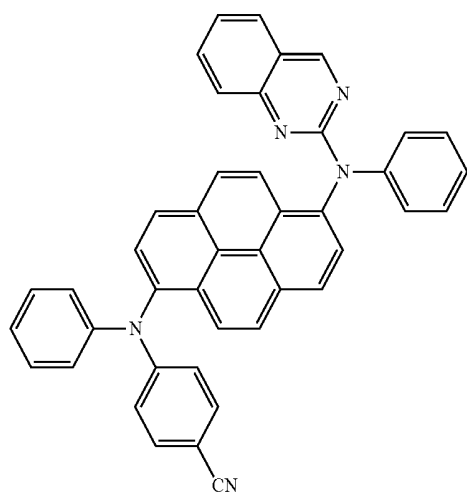


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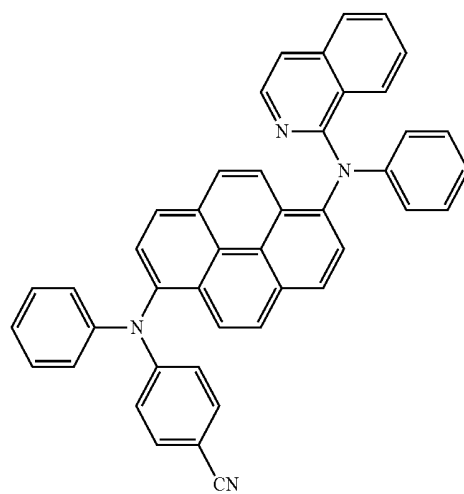
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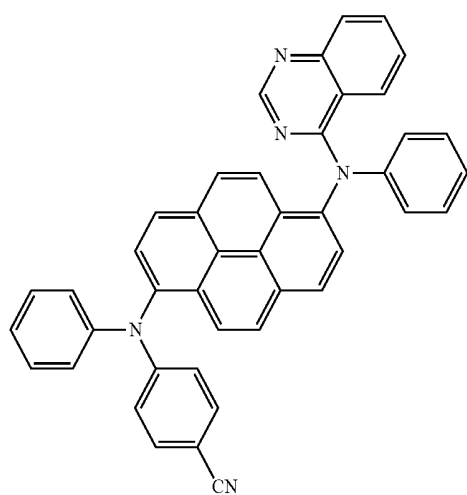
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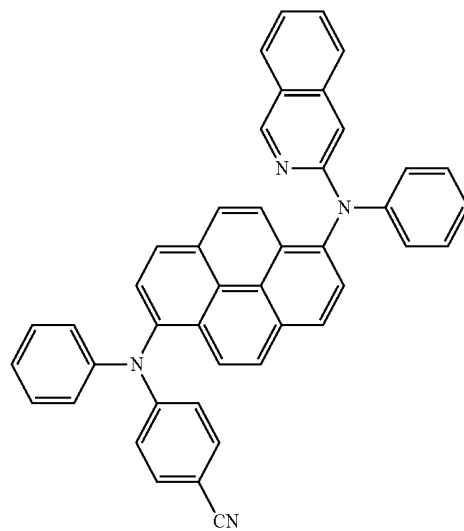
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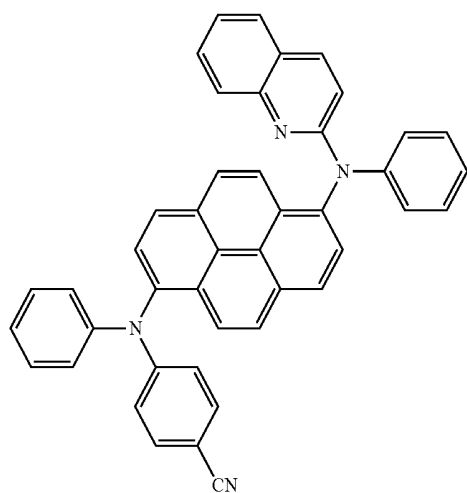


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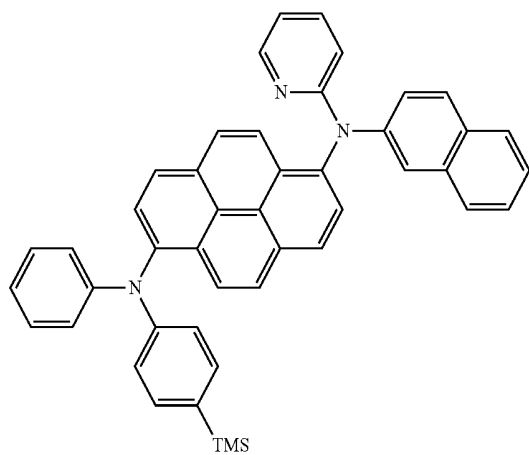


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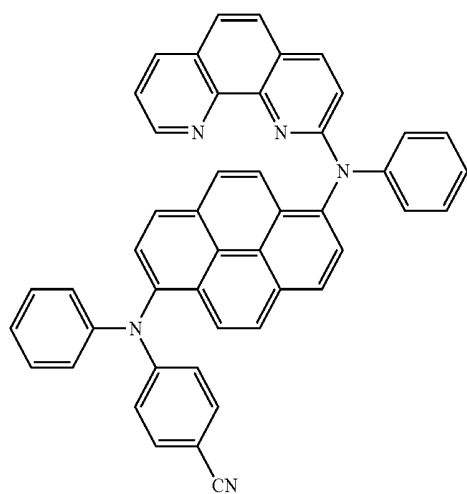
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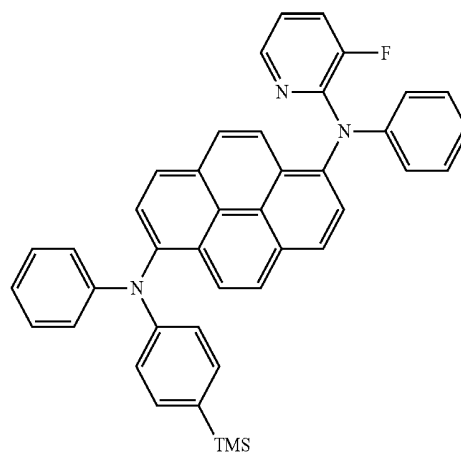
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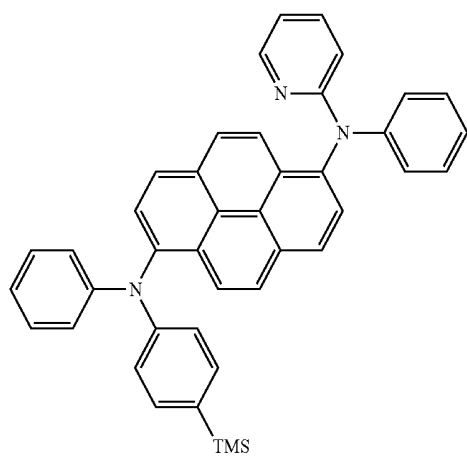
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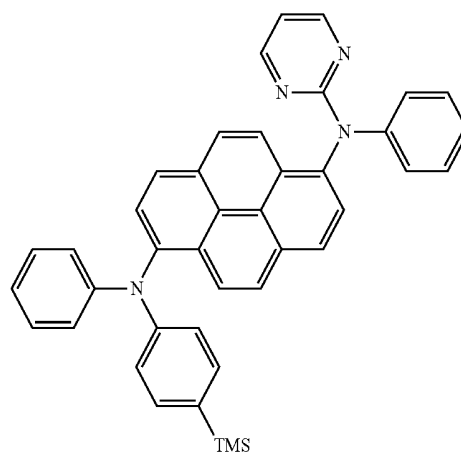
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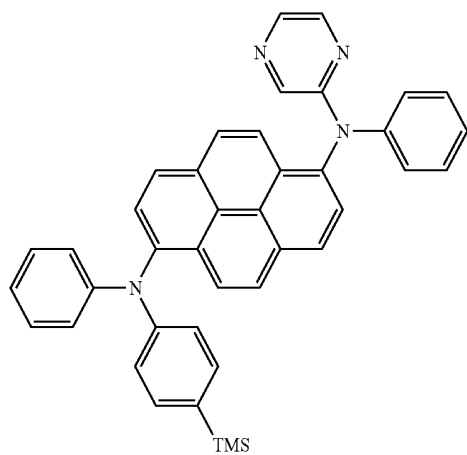
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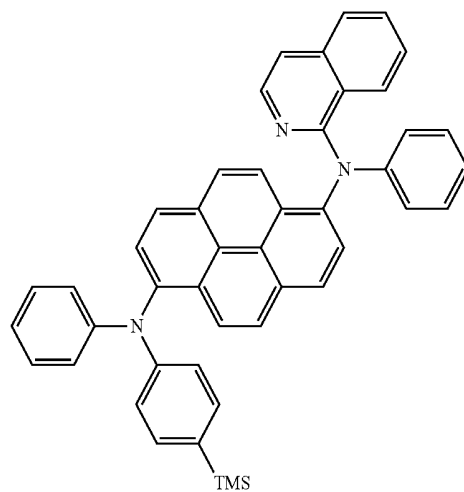
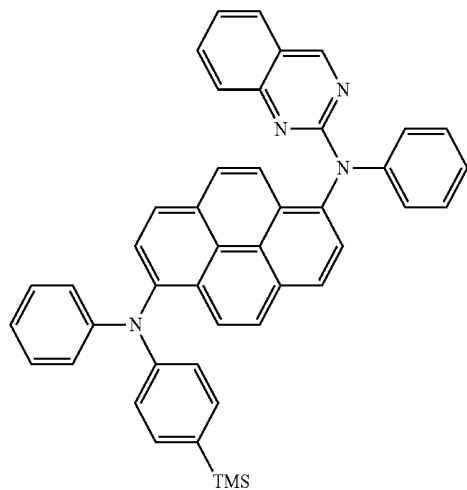
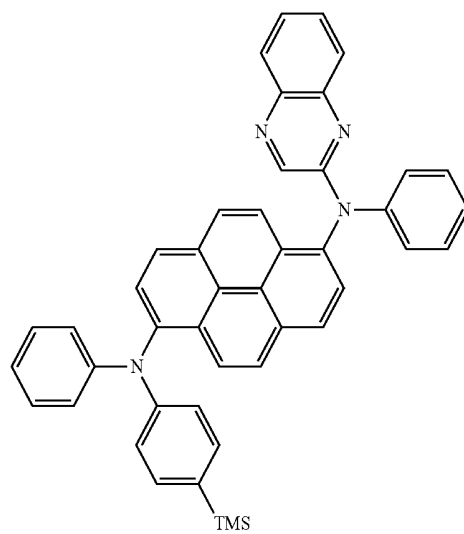
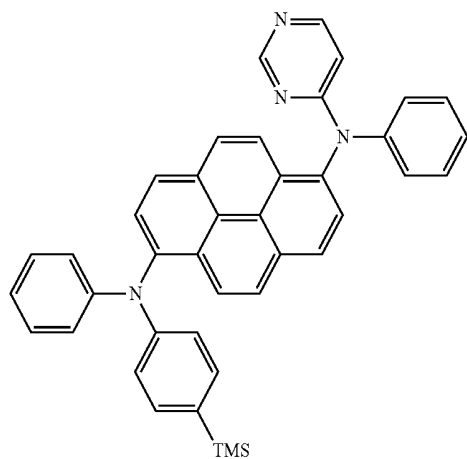
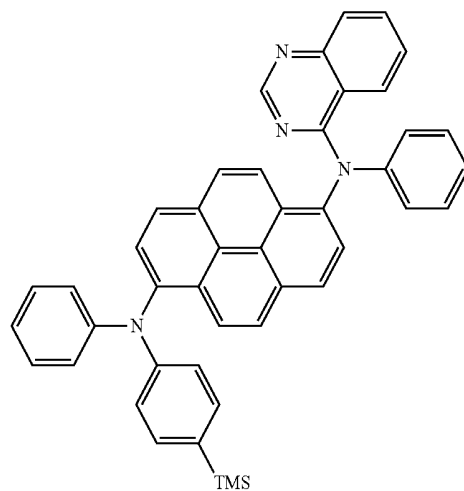
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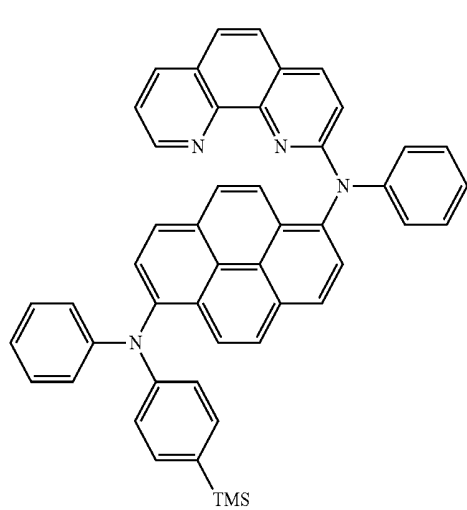
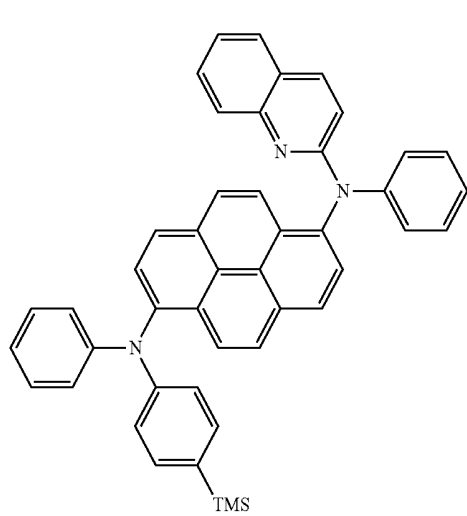
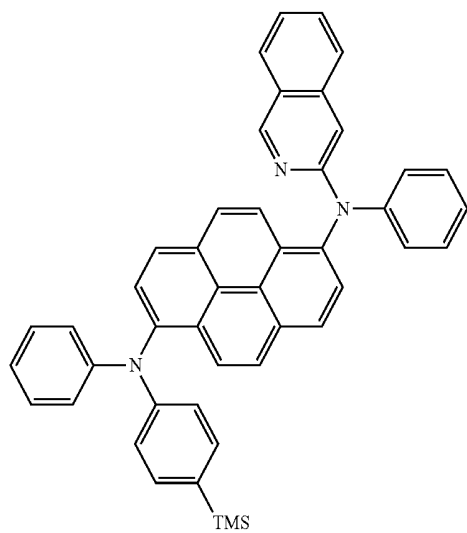
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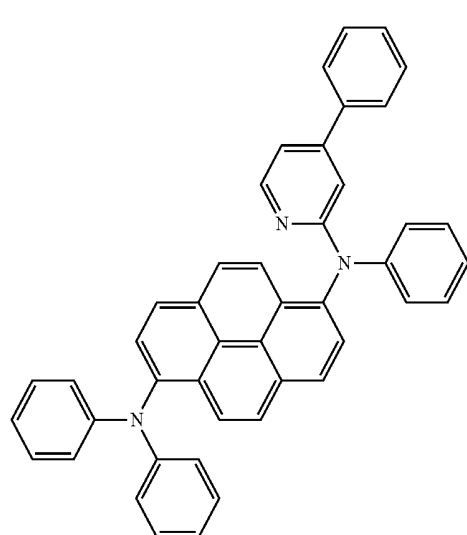
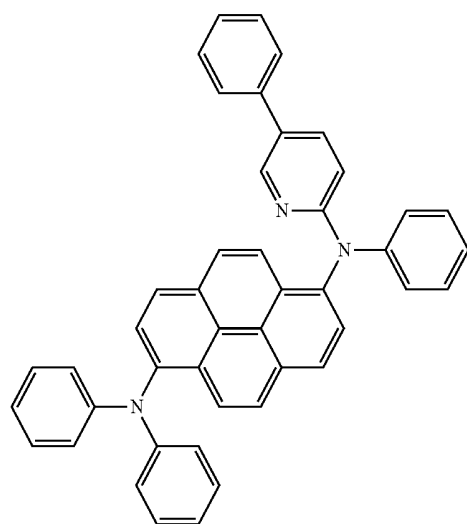
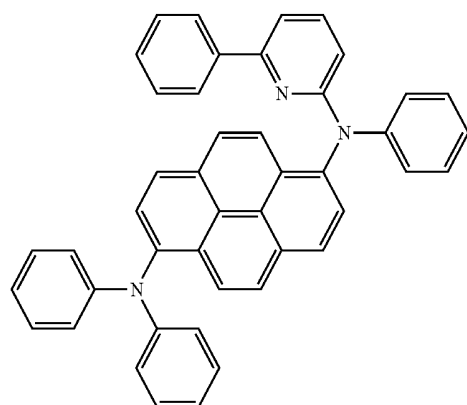
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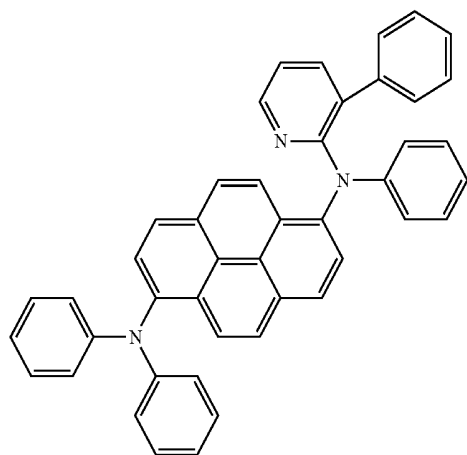
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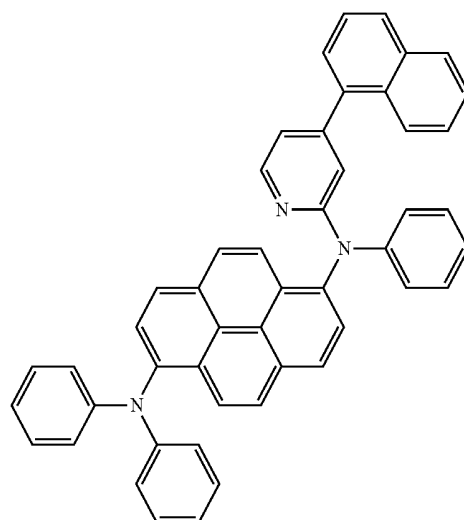


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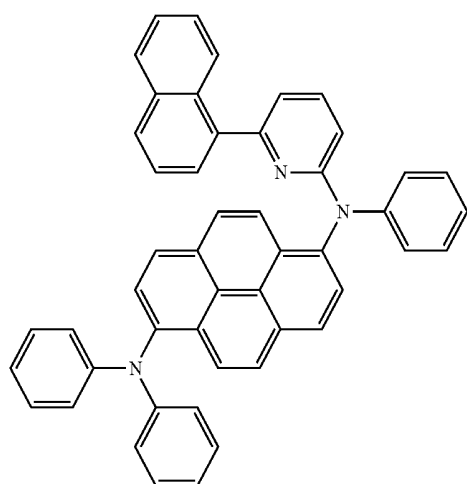


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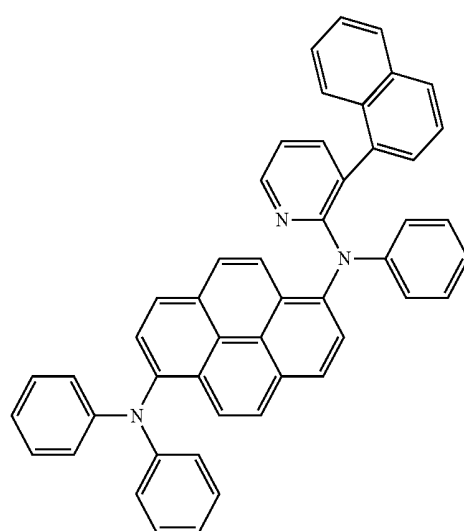
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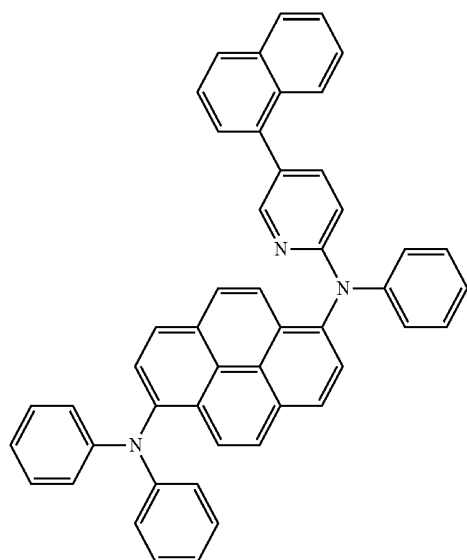
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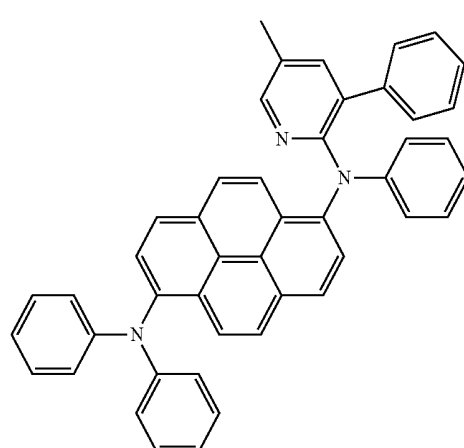
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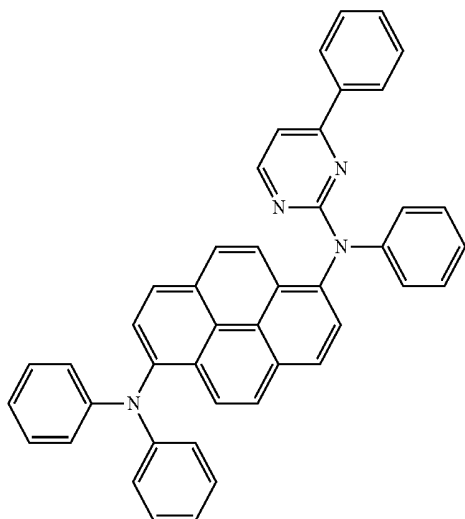


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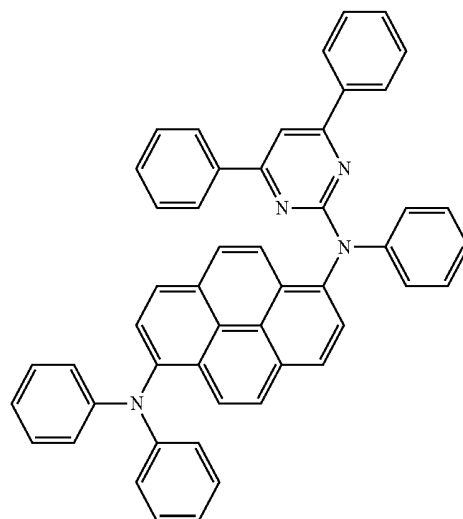
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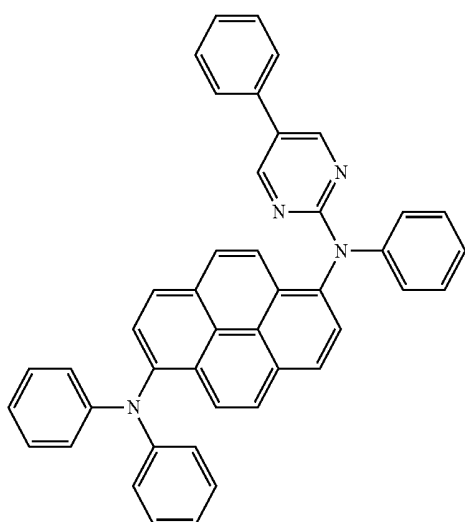


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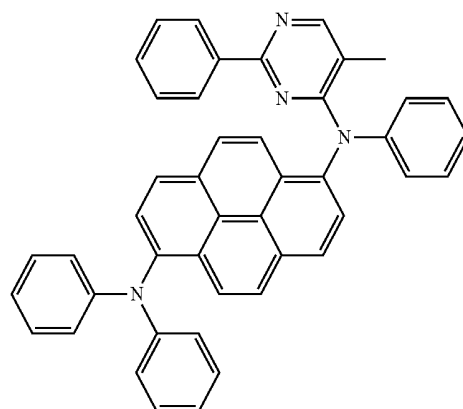
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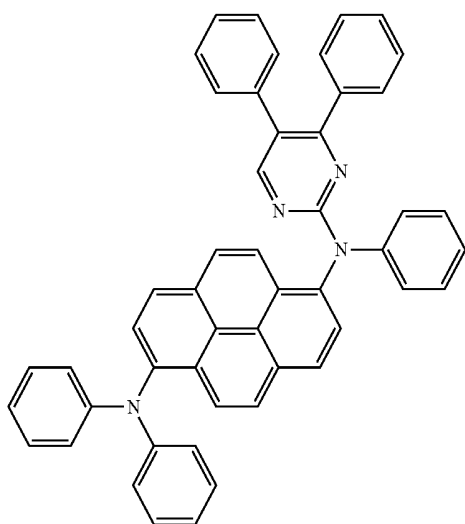
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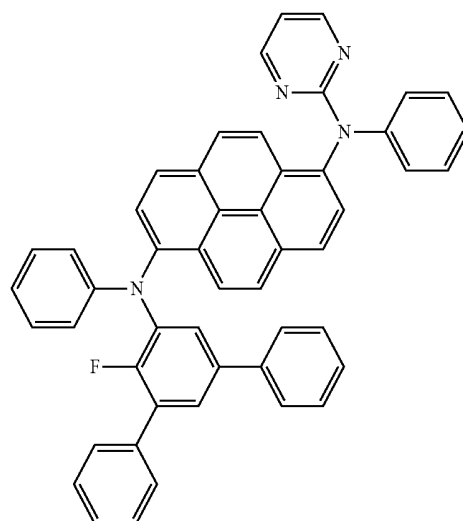
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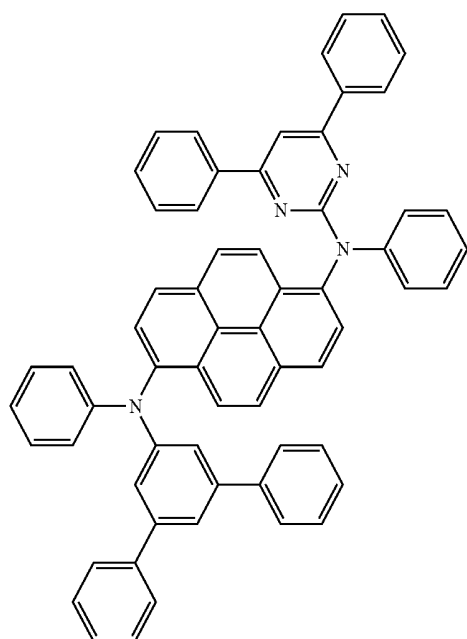
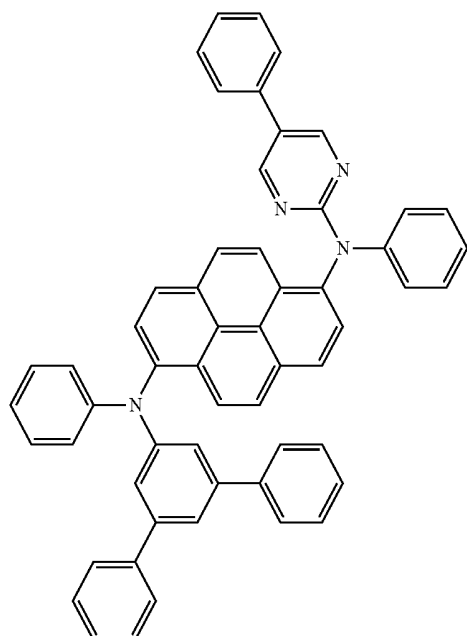


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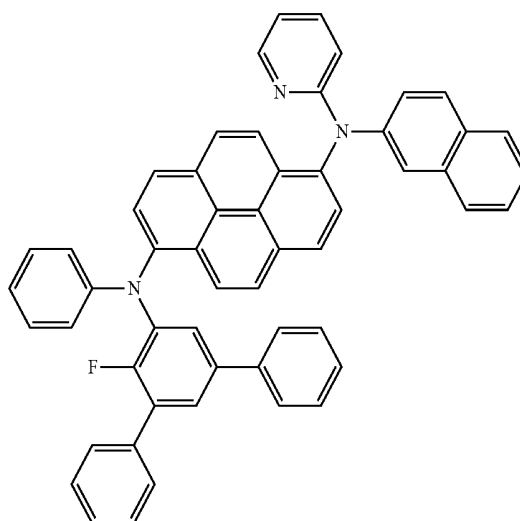
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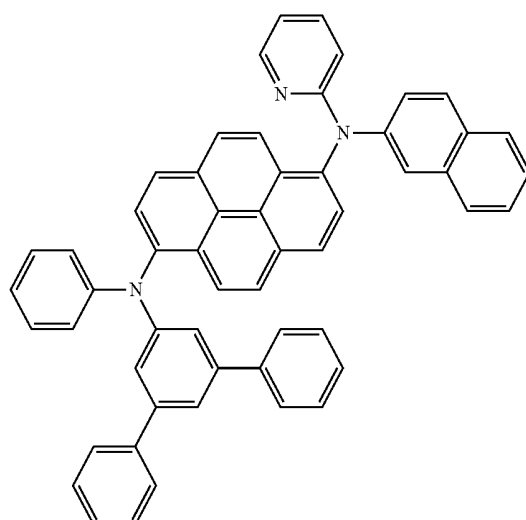


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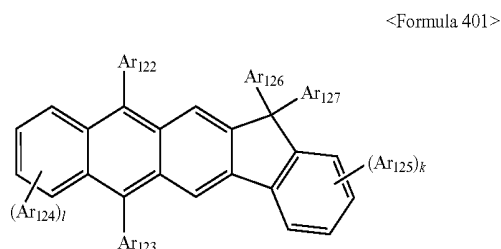
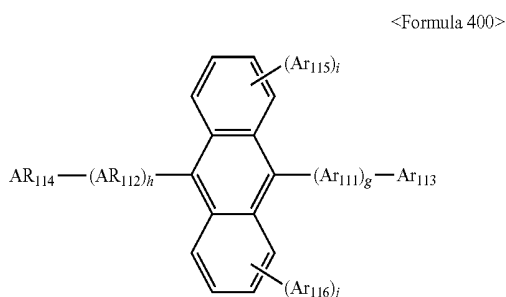


[0027] The embodiments may also be realized by providing an organic light-emitting diode including a first electrode; a second electrode facing the first electrode; and an organic layer between the first electrode and the second electrode and including an emission layer, wherein the organic layer includes the pyrene-based compound according to an embodiment.

[0028] The organic layer may further include a hole transportation region between the first electrode and the emission layer, the hole transportation region including at least one of a hole injection layer, a hole transportation layer, a functional layer having a hole injection capability and a hole transport capability, a buffer layer, and an electron blocking layer, and an electron transportation region between the emission layer and the second electrode, the electron transportation region including at least one of a hole blocking layer, an electron transportation layer, and an electron injection layer.

[0029] The pyrene-based compound may be in the emission layer, and a fluorescent dopant that emits light according to a fluorescence emission mechanism, the emission layer further including a host.

[0030] The host may include at least one of an anthracene-based compound represented by Formula 400, below, or an anthracene-based compound represented by Formula 401, below:



[0031] wherein, in Formulae 400 and 401, Ar_{111} 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 l are each independently an integer of 0 to 4.

BRIEF DESCRIPTION OF THE DRAWING

[0032] Features will be apparent to those of skill in the art by describing in detail exemplary embodiments with reference to the attached drawing in which:

[0033] FIG. 1 illustrates a cross-sectional view of the structure of an organic light-emitting diode according to an embodiment.

DETAILED DESCRIPTION

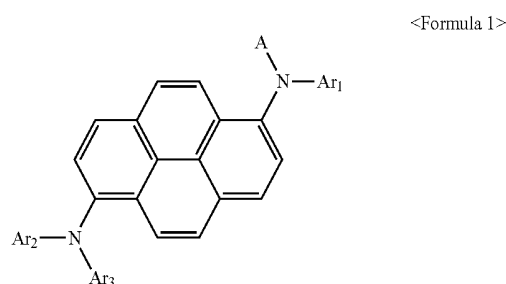
[0034] Example embodiments will now be described more fully hereinafter with reference to the accompanying drawing; however, they may be embodied in different forms and should not be construed as limited to the embodiments set forth herein. Rather, these embodiments are provided so that this disclosure will be thorough and complete, and will fully convey exemplary implementations to those skilled in the art.

[0035] In the drawing figures, the dimensions of layers and regions may be exaggerated for clarity of illustration. It will also be understood that when a layer or element is referred to as being “on” another layer or substrate, it can be directly on the other layer or substrate, or intervening layers may also be present. Further, it will be understood that when a layer is referred to as being “under” another layer, it can be directly under, and one or more intervening layers may also be present. In addition, it will also be understood that when a layer is referred to as being “between” two layers, it can be the

only layer between the two layers, or one or more intervening layers may also be present. Like reference numerals refer to like elements throughout.

[0036] As used herein, the term “and/or” includes any and all combinations of one or more of the associated listed items. Expressions such as “at least one of,” when preceding a list of elements, modify the entire list of elements and do not modify the individual elements of the list.

[0037] A pyrene-based compound according to an embodiment may be represented by Formula 1, below.



[0038] In Formula 1, ring A may be a substituted or unsubstituted C_2-C_{30} heteroaromatic group including at least one nitrogen (N).

[0039] In Formula 1-1, Ar_{1n} in an implementation, A may be selected from a substituted or unsubstituted pyridinyl group, a substituted or unsubstituted pyrimidinyl group, a substituted or unsubstituted triazinyl group, a substituted or unsubstituted pyrazinyl group, a substituted or unsubstituted pyridazinyl group, a substituted or unsubstituted isindolyl group, a substituted or unsubstituted indolyl group, a substituted or unsubstituted indazolyl group, a substituted or unsubstituted isoquinolinyl group, a substituted or unsubstituted quinolinyl group, a substituted or unsubstituted phthalazinyl group, a substituted or unsubstituted quinoxalinyl group, a substituted or unsubstituted quinazolinyl group, a substituted or unsubstituted cinnolinyl group, or a substituted or unsubstituted 1,10-phenanthrolinyl.

[0040] For example, A may be selected from:

[0041] a pyridinyl group, a pyrimidinyl group, a triazinyl group, a pyrazinyl group, a pyridazinyl group, an isindolyl group, an indolyl group, an indazolyl group, an isoquinolinyl group, a quinolinyl group, a phthalazinyl group, a quinoxalinyl group, a cinnolinyl group, and a 1,10-phenanthrolinyl group; and

[0042] a pyridinyl group, a pyrimidinyl group, a triazinyl group, a pyrazinyl group, a pyridazinyl group, an isindolyl group, an indolyl group, an indazolyl group, an isoquinolinyl group, a quinolinyl group, a phthalazinyl group, a quinoxalinyl group, a cinnolinyl group, and a 1,10-phenanthrolinyl group, each substituted with at least one of:

[0043] a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C_1-C_{60} alkyl group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, and a C_1-C_{60} alkoxy group,

[0044] a C_1-C_{60} alkyl group, a C_2-C_{60} alkenyl group, a C_2-C_{60} alkynyl group, and a C_1-C_{60} alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone

group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof,

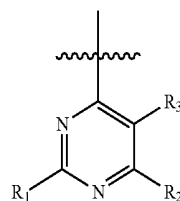
[0045] a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, and a C₂-C₆₀ heteroaryl group,

[0046] a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, and a C₂-C₆₀ heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group, and

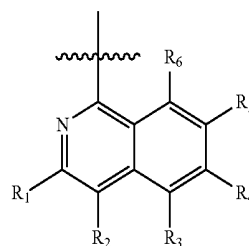
[0047] —Si(Q₁₁)(Q₁₂)(Q₁₃) (in which Q₁₁ to Q₁₃ are each independently a C₁-C₆₀ alkyl group, a C₁-C₆₀ alkoxy group, a C₆-C₆₀ aryl group, or a C₂-C₆₀ heteroaryl group).

[0048] In an implementation, A in Formula 1 may be represented by one of Formulae 2A to 2K below, in which the broken line represents a bonding location with Formula 1.

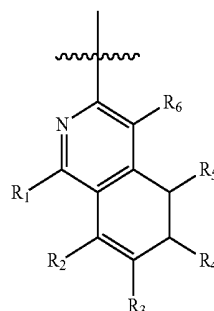
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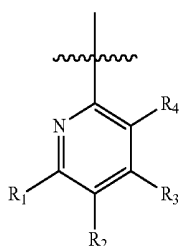
Formula 2D



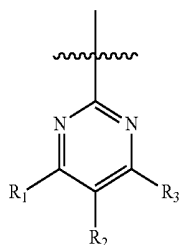
Formula 2E



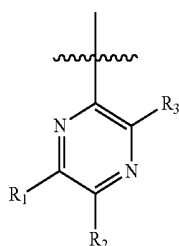
Formula 2F



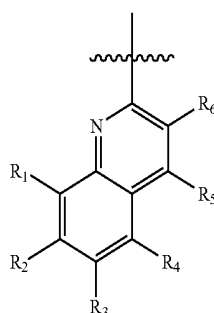
Formula 2A



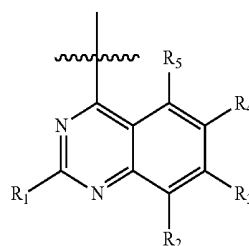
Formula 2B



Formula 2C

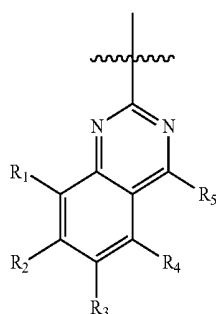


Formula 2G

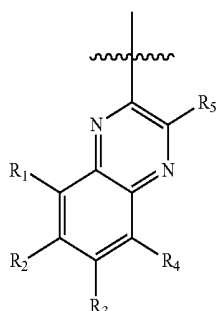


Formula 2H

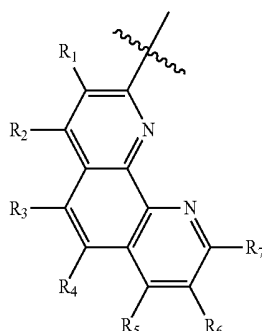
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Formula 2I



Formula 2J



Formula 2K

[0049] R_1 to R_7 in Formulae 2A to 2K may each independently be selected from:

[0050] a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group;

[0051] a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof;

[0052] a C_3 - C_{10} cycloalkyl group, a C_3 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_3 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, and a C_2 - C_{60} heteroaryl group;

[0053] a C_3 - C_{10} cycloalkyl group, a C_3 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_3 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, and a C_2 - C_{60} heteroaryl group, each substituted with at least one of a deuterium atom,

a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic 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, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group; and

[0054] $-\text{Si}(\text{Q}_{11})(\text{Q}_{12})(\text{Q}_{13})$ (in which Q_{11} to Q_{13} are each independently a C_1 - C_{60} alkyl group, a C_1 - C_{60} alkoxy group, a C_6 - C_{60} aryl group, or a C_2 - C_{60} heteroaryl group).

[0055] In an implementation, R_1 to R_7 in Formulae 2A to 2K may each independently be selected from:

[0056] a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a C_1 - C_{10} alkyl group;

[0057] a C_1 - C_{60} alkyl group substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof;

[0058] a C_6 - C_{16} aryl group and a C_2 - C_{16} heteroaryl group;

[0059] a C_6 - C_{16} aryl group and a C_2 - C_{16} heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C_1 - C_{10} alkyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group; and

[0060] $-\text{Si}(\text{Q}_{11})(\text{Q}_{12})(\text{Q}_{13})$ (in which Q_{11} to Q_{13} are each independently a C_1 - C_{10} alkyl group, a C_6 - C_{16} aryl group, or a C_2 - C_{16} heteroaryl group), but are not limited thereto.

[0061] For example, R_1 to R_7 in Formulae 2A to 2K may each independently be selected from:

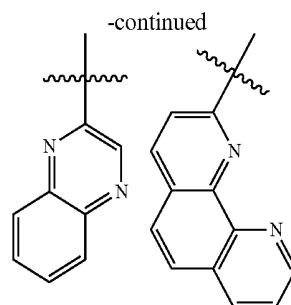
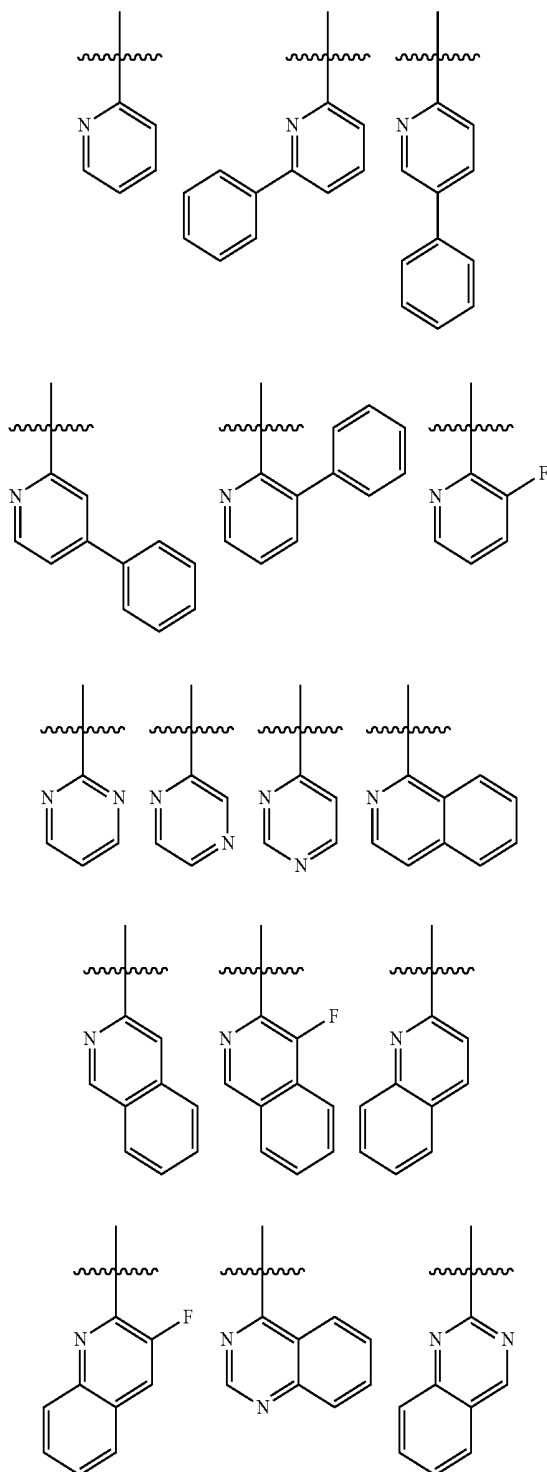
[0062] a hydrogen atom, a deuterium atom, F, Cl, Br, I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof, a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, and a hexyl group;

[0063] a phenyl group, a naphthyl group, and an anthracenyl group; and

[0064] a phenyl group, a naphthyl group, and an anthracenyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, a hexyl group, phenyl group, a naphthyl group, an anthracenyl group,

a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, and a phenylcarbazolyl group.

[0065] According to an embodiment, A in Formula 1 may have one of the following structures, in which the broken line represents a bonding location with Formula 1.



[0066] Ar₁, Ar₂, and Ar₃ in Formula 1 may each independently be selected from a substituted or unsubstituted phenyl group and a substituted or unsubstituted naphthyl group.

[0067] For example, Ar₁, Ar₂ and Ar₃ in Formula 1 may each independently be selected from:

[0068] a phenyl group and a naphthyl group; and

[0069] a phenyl group and a naphthyl group, each substituted with at least one of:

[0070] a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group,

[0071] a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof,

[0072] a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, and a C₂-C₆₀ heteroaryl group,

[0073] a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, and a C₂-C₆₀ heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group, and

[0074] —Si(Q₁₁)(Q₁₂)(Q₁₃) (in which Q₁₁ to Q₁₃ are each independently a C₁-C₆₀ alkyl group, a C₁-C₆₀ alkoxy group, a C₆-C₆₀ aryl group, or a C₂-C₆₀ heteroaryl group).

[0075] For example, Ar₁, Ar₂ and Ar₃ in Formula 1 may each independently be selected from:

[0076] a phenyl group and a naphthyl group; and

[0077] a phenyl group and a naphthyl group, each substituted with at least one of:

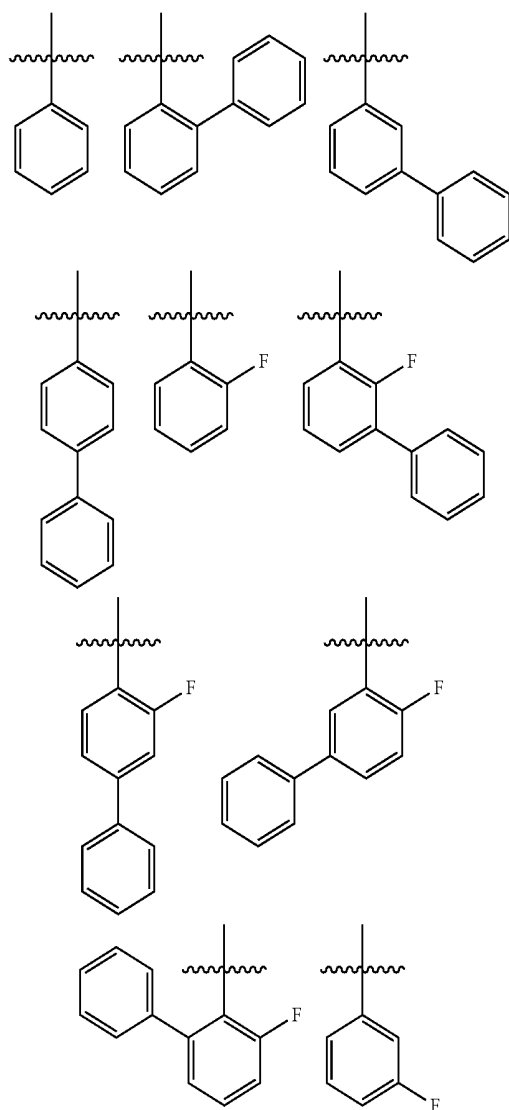
[0078] a deuterium atom, F, Cl, Br, I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino

group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof and a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, and a hexyl group,

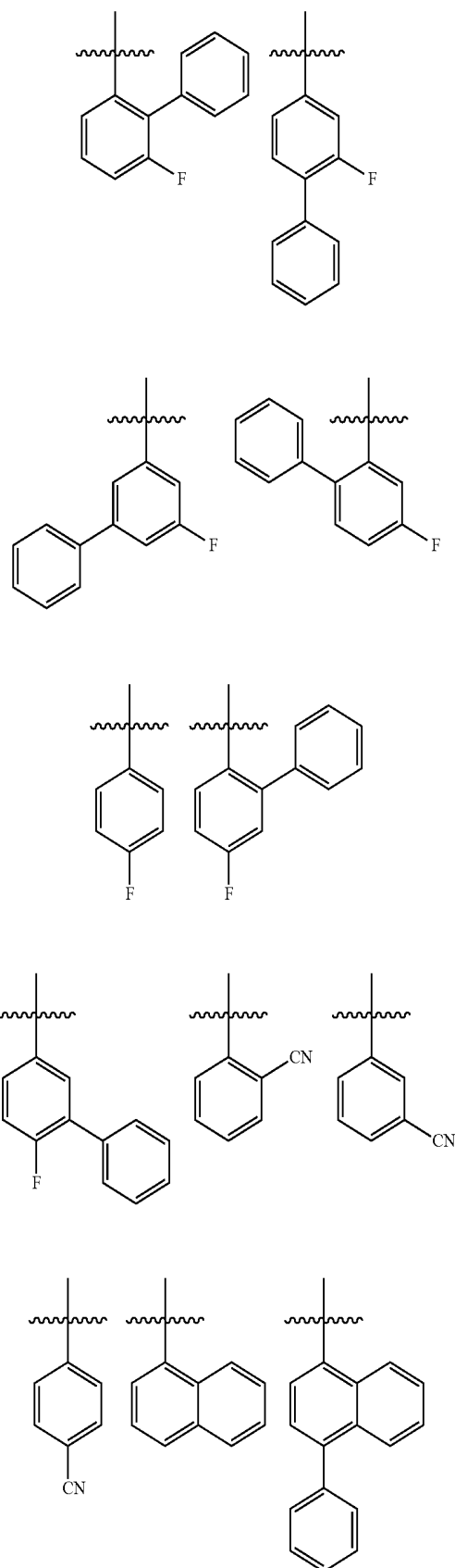
[0079] a phenyl group, a naphthyl group, and an anthracenyl group, and

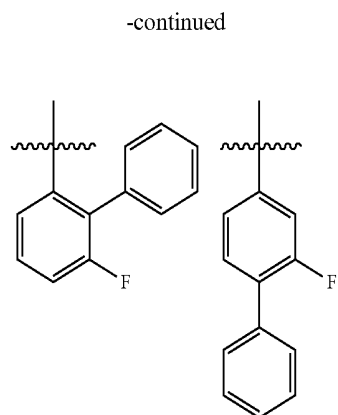
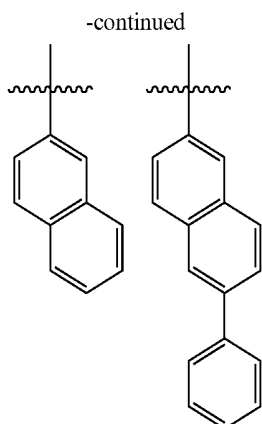
[0080] a phenyl group, a naphthyl group, and an anthracenyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, a hexyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, and a phenylcarbazolyl group.

[0081] In an implementation, Ar₁ in Formula 1 may have one of the following structures, in which the broken line represents a bonding location with Formula 1.

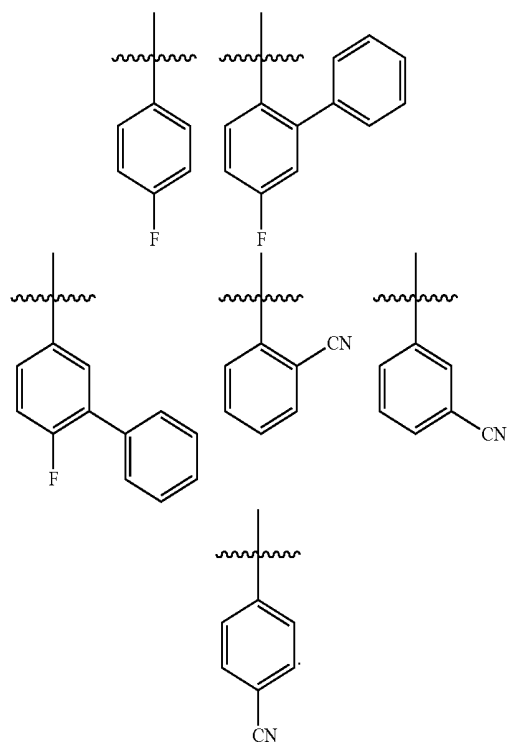
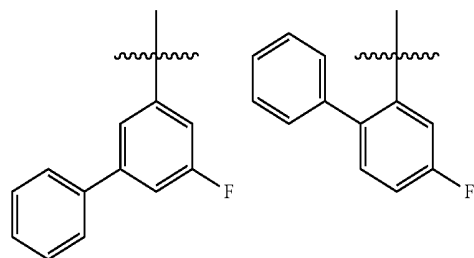
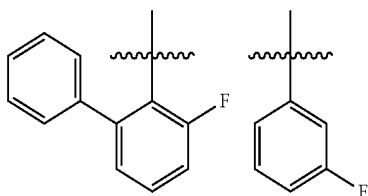
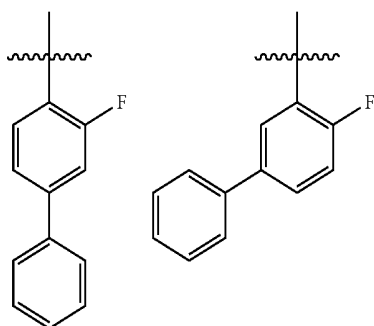
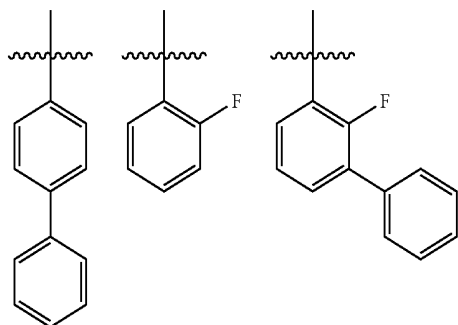
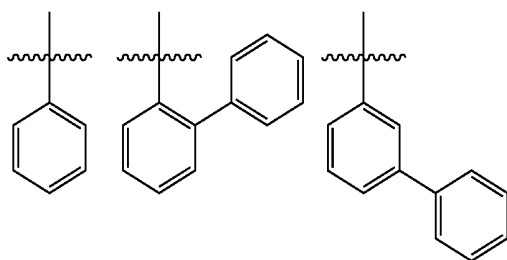


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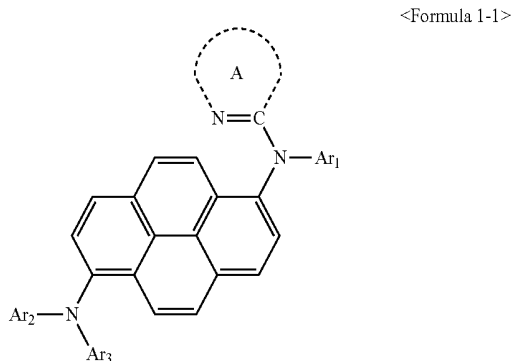


[0082] In an implementation, Ar_2 and Ar_3 in Formula 1 may each independently have one of the following structures, in which the broken line represents a bonding location with Formula 1.



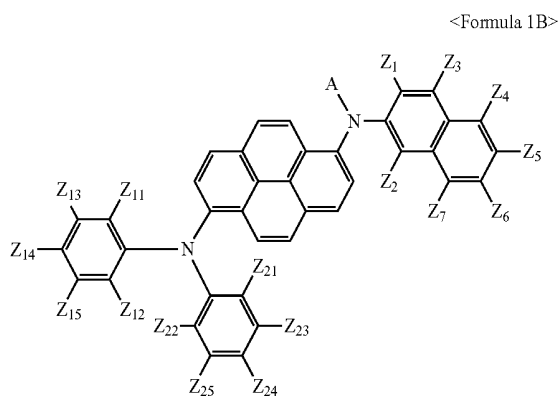
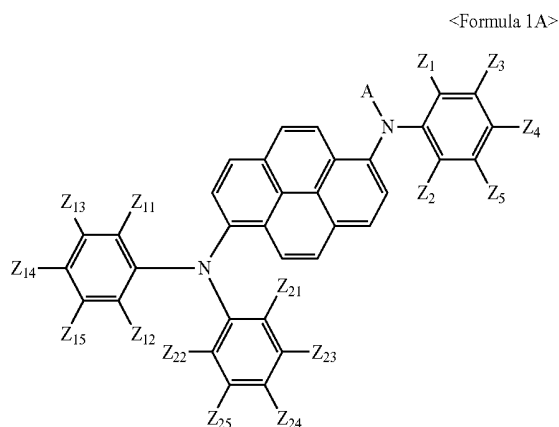
[0083] In Formula 1, Ar_1 may be a substituted or unsubstituted phenyl group or a substituted or unsubstituted naphthyl group, and Ar_2 and Ar_3 may each independently be a substituted or unsubstituted phenyl group.

[0084] In an implementation, the pyrene-based compound represented by Formula 1 may be represented by Formula 1-1, below.

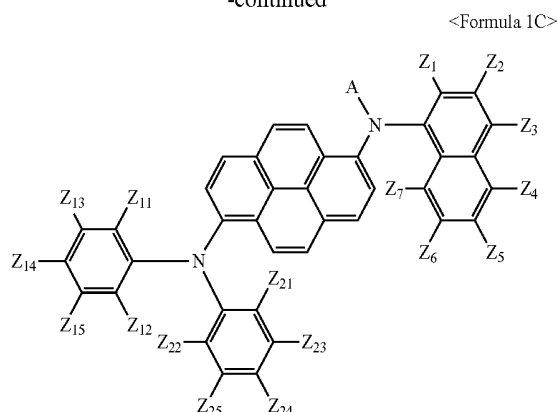


[0085] In Formula 1-1, Ar₁, Ar₂, and Ar₃ may be the same as described with respect to Formula 1, and A is shown as a partially formed heterocyclic ring.

[0086] According to an embodiment, the pyrene-based compound may be represented by one of Formulae 1A, 1B, or 1C, below.



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[0087] A in Formulae 1A, 1B, and 1C may be as described above with respect to Formula 1. For example, A in Formulae 1A, 1B, and 1C may be represented by one of Formulae 2A to 2K, above.

[0088] Z₁ to Z₇, Z₁₁ to Z₁₅, and Z₂₁ to Z₂₅ in Formulae 1A, 1B, and 1C may each independently be selected from a pyridinyl group, a pyrimidinyl group, a triazinyl group, a pyrazinyl group, a pyridazinyl group, an isoindolyl group, an indolyl group, an indazolyl group, an isoquinolinyl group, a quinolinyl group, a phthalazinyl group, a quinoxalinyl group, a cinnolinyl group, and a 1,10-phenanthrolinyl group. Each may be substituted with at least one of i) a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine, a hydrazone, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₆₀ alkoxy group; ii) a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof; iii) a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group and a C₂-C₆₀ heteroaryl group; iv) a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group and a C₂-C₆₀ heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group; and v) —Si(Q₁₁)(Q₁₂)(Q₁₃) (in which Q₁₁ to Q₁₃ are each independently a C₁-C₆₀ alkyl group, a C₁-C₆₀ alkoxy group, a C₆-C₆₀ aryl group, or a C₂-C₆₀ heteroaryl group).

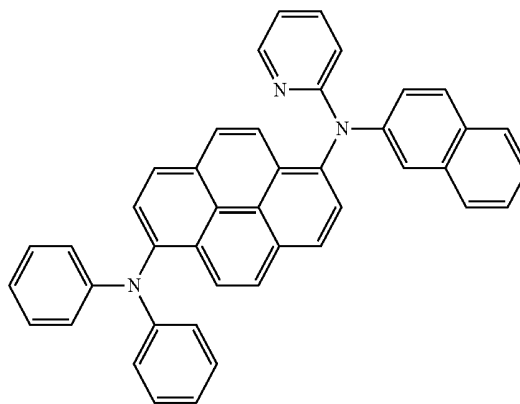
[0089] For example, Z_1 to Z_7 , Z_{11} to Z_{15} , and Z_{21} to Z_{25} in Formulae 1A, 1B and 1C may each independently be selected from a phenyl group and a naphthyl group. Each may substituted with at least one of i) a hydrogen atom, a deuterium atom, F, Cl, Br, I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a phosphoric acid or a salt thereof and a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, and a hexyl group; ii) a phenyl group, a naphthyl group, and an anthracenyl group; and iii) a phenyl group, a naphthyl group and an anthracenyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, a hexyl group, a phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, and a phenylcarbazolyl group.

[0090] According to an embodiment, the pyrene-based compound may be represented by Formula 1A, 1B, or 1C, A may be represented by one of Formulae 2A to 2K, and R_1 to R_6 , Z_1 to Z_7 , Z_{11} to Z_{15} and Z_{21} to Z_{25} may each independently be a hydrogen atom, a F atom, a cyano group, or a phenyl group in the Formula 1A, 1B, and 1C.

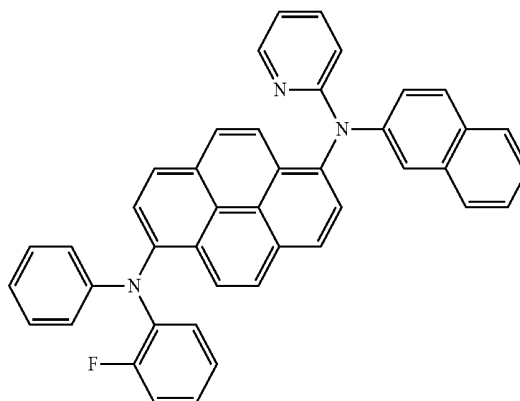
[0091] According to another embodiment, the pyrene-based compound may be one of Compounds 1 to 96 illustrated below.

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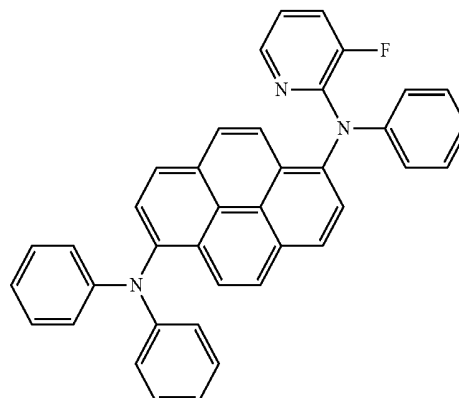
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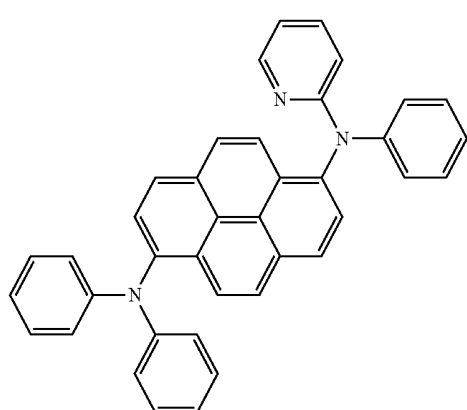
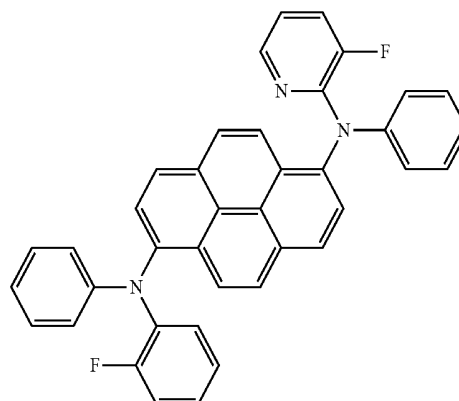
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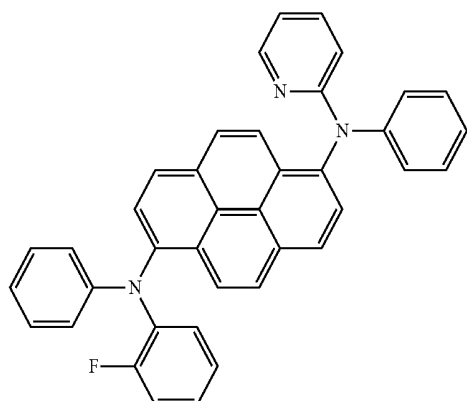
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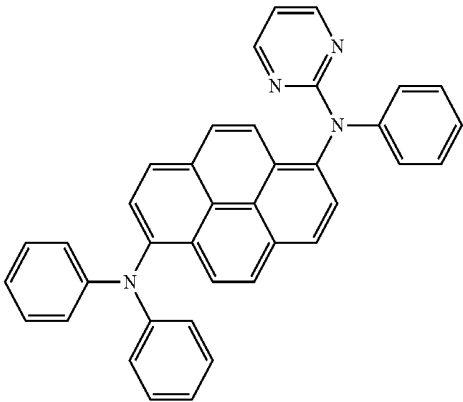


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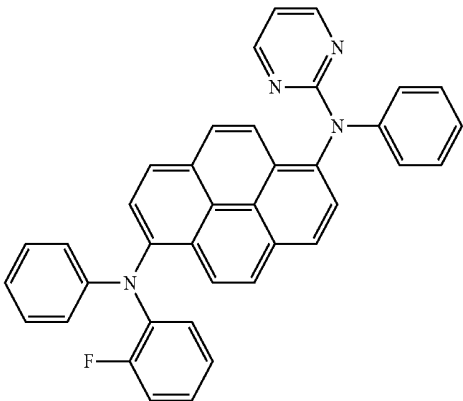


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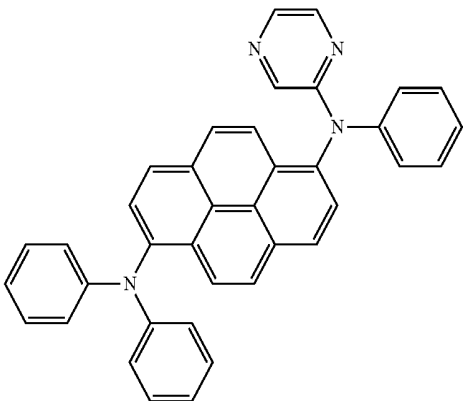
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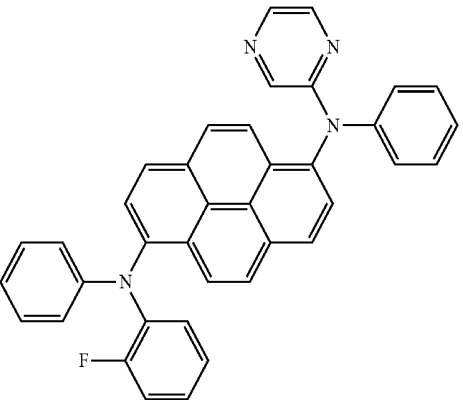
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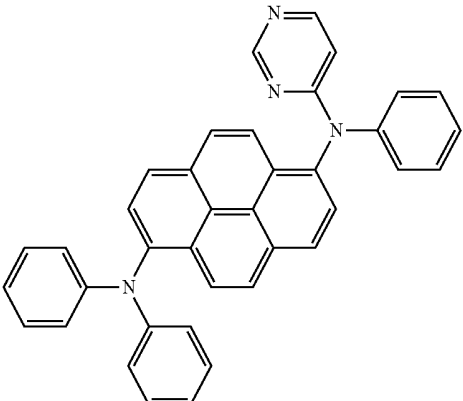


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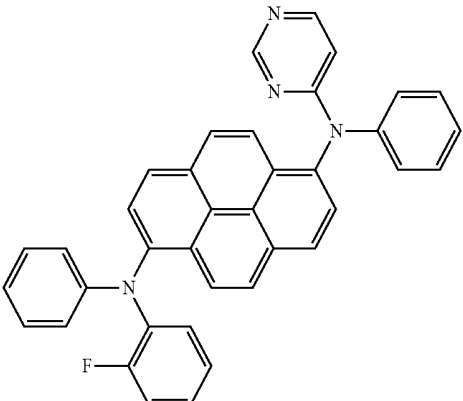


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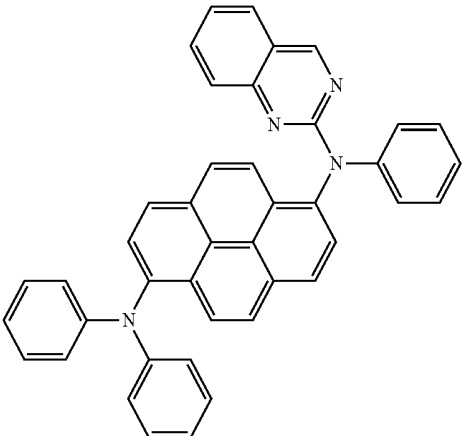
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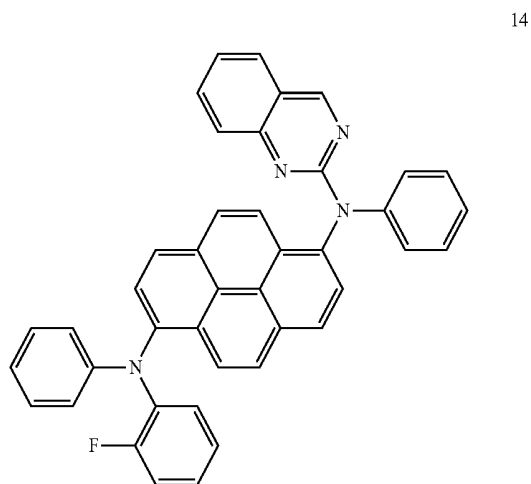


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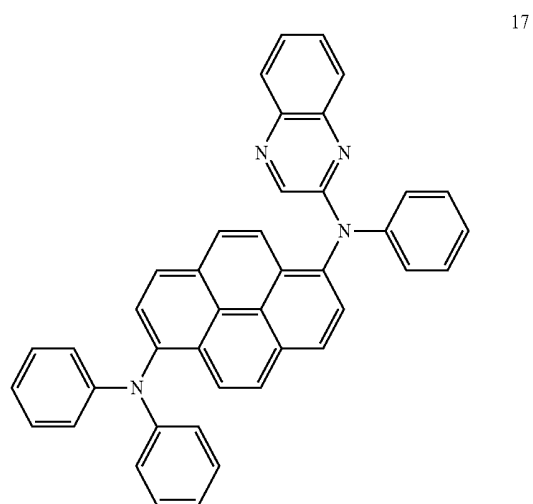


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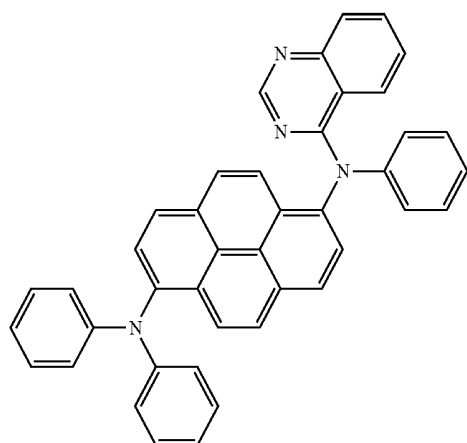
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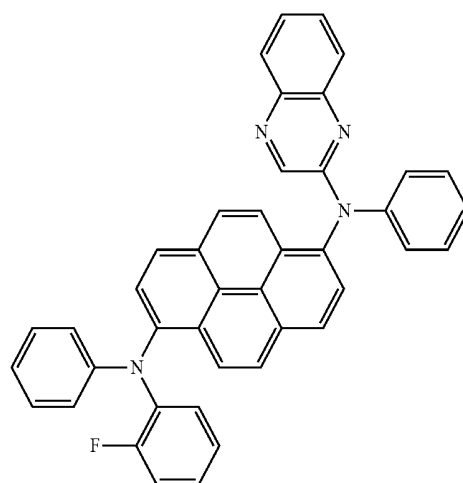
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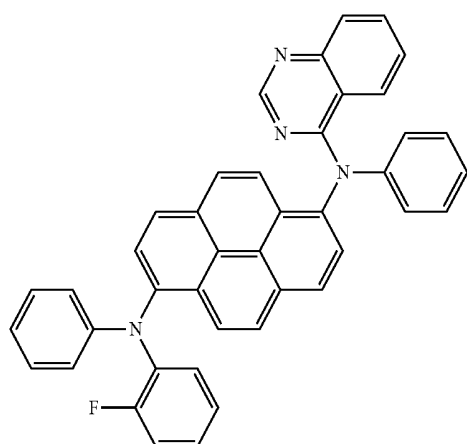
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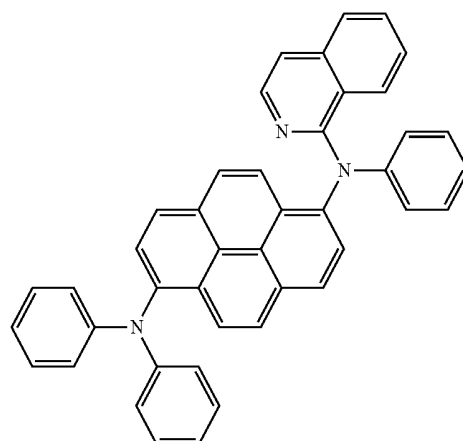
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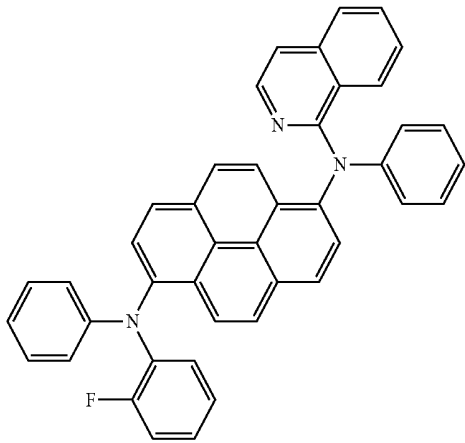


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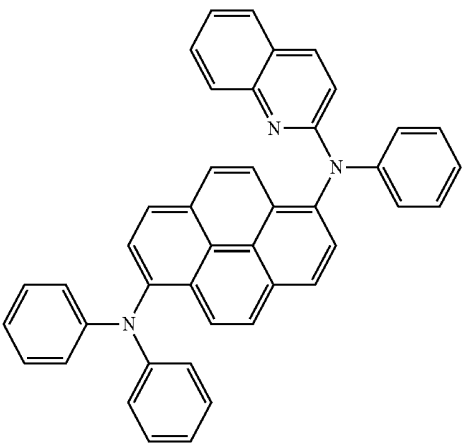


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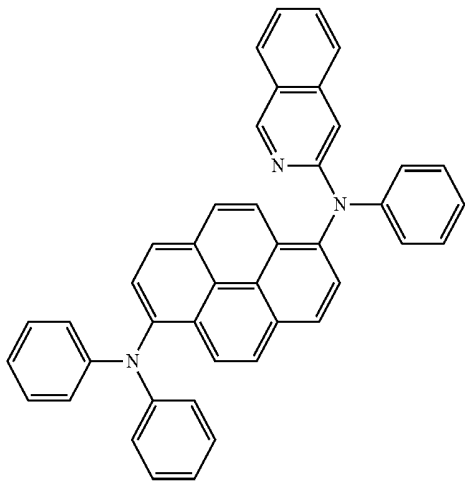
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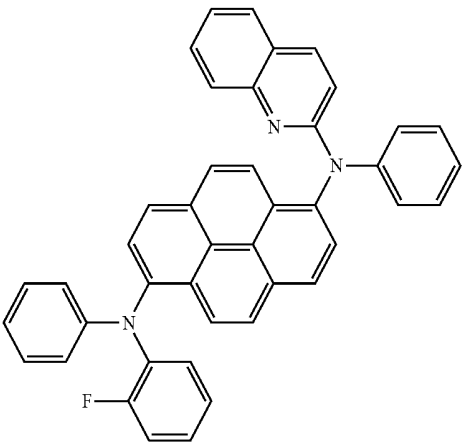
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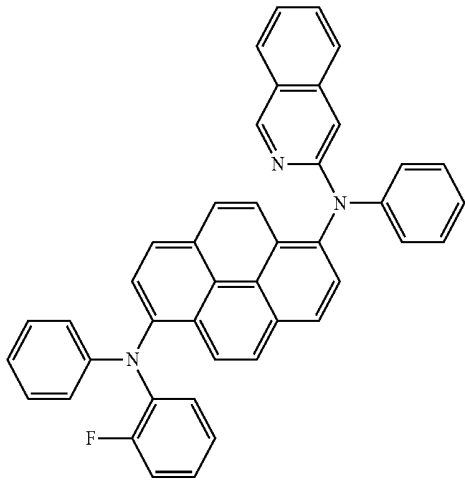
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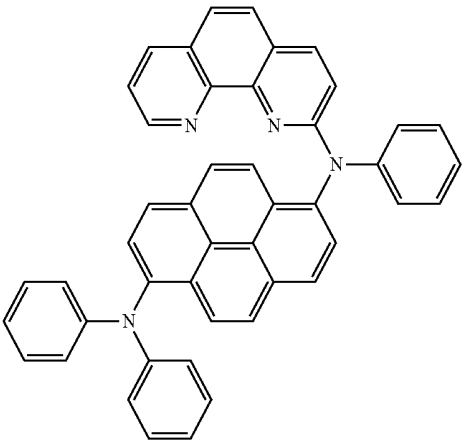
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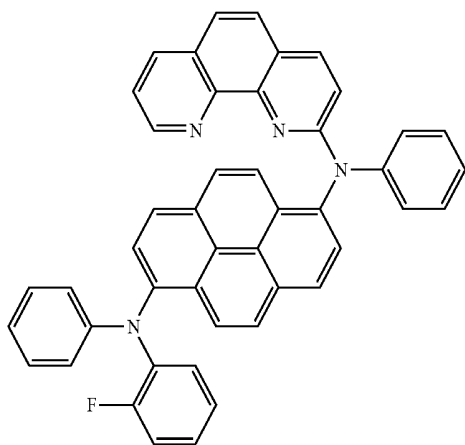


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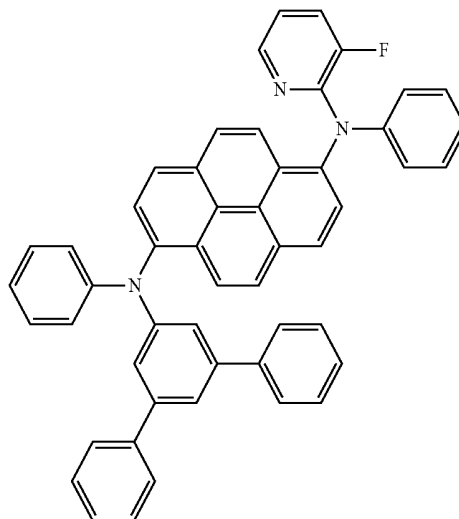
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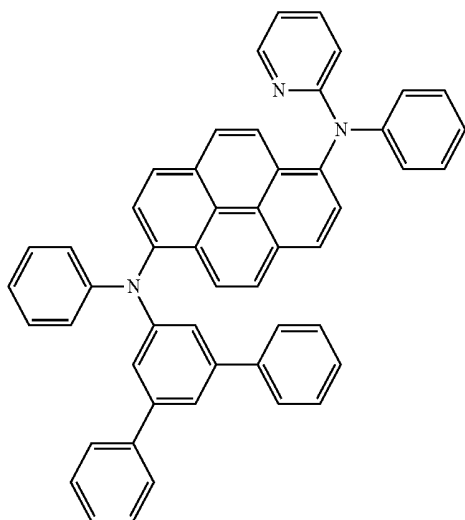


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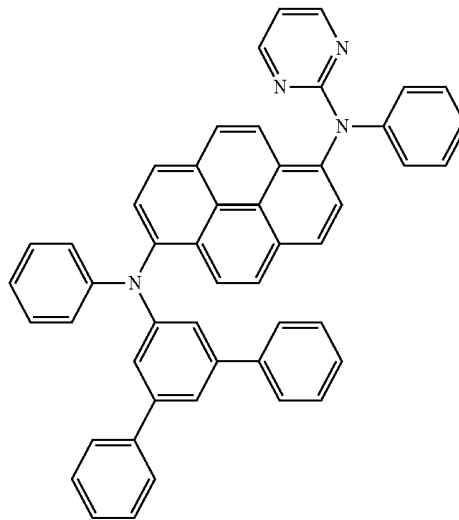
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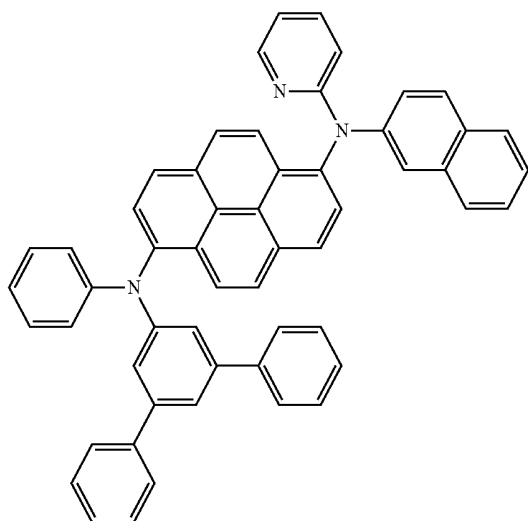
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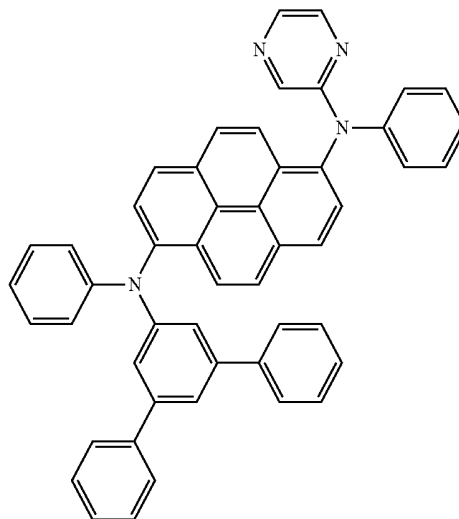
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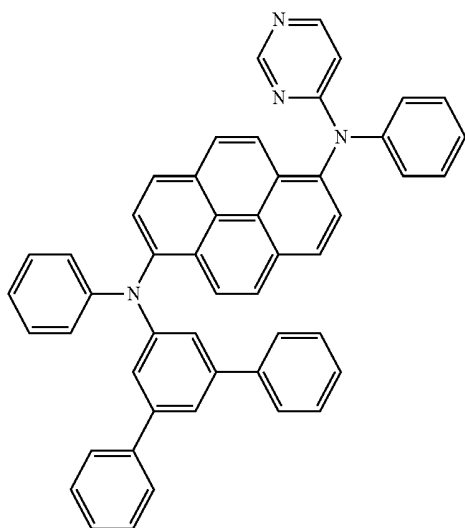
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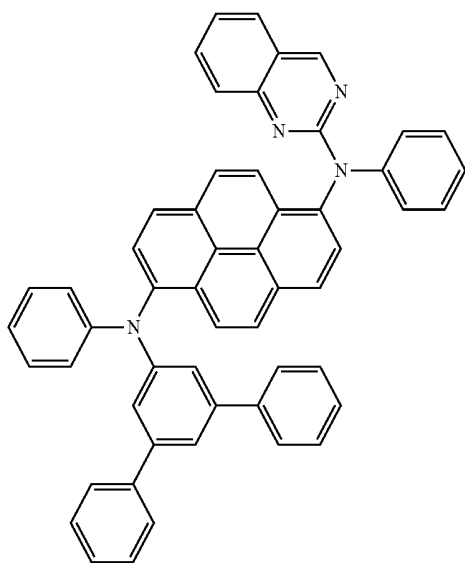
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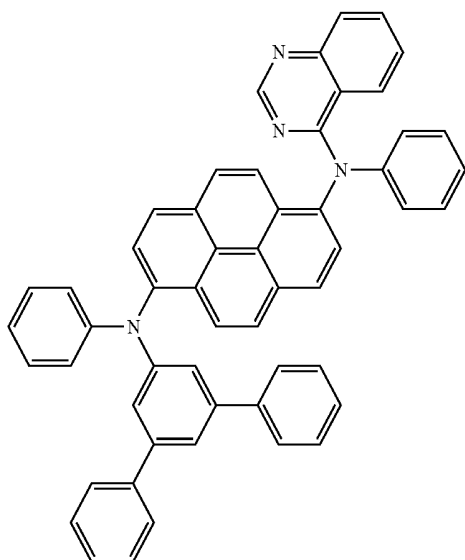
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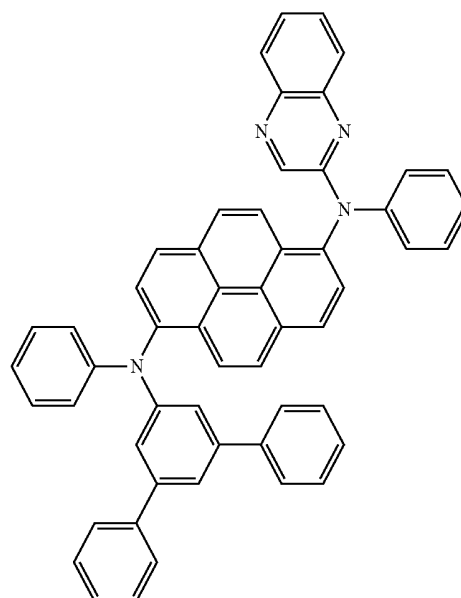


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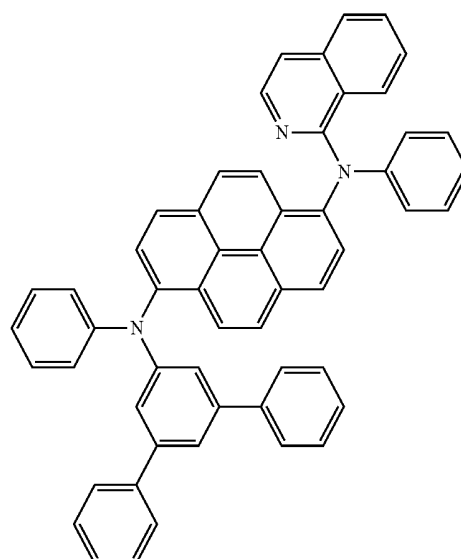


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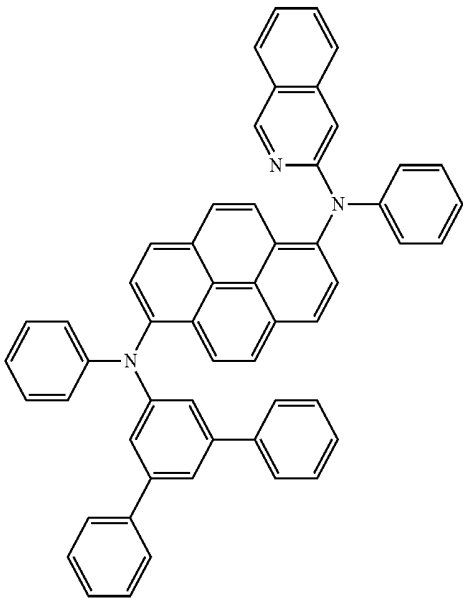
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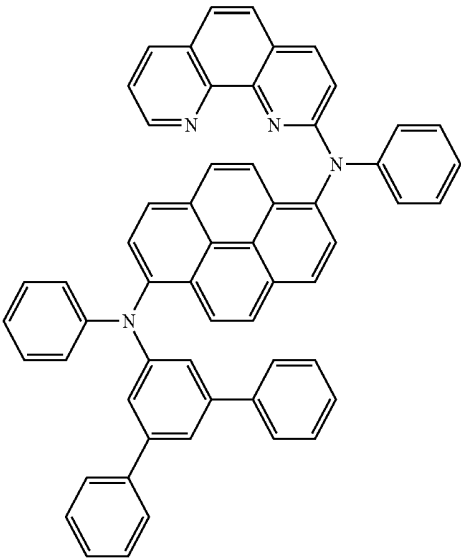
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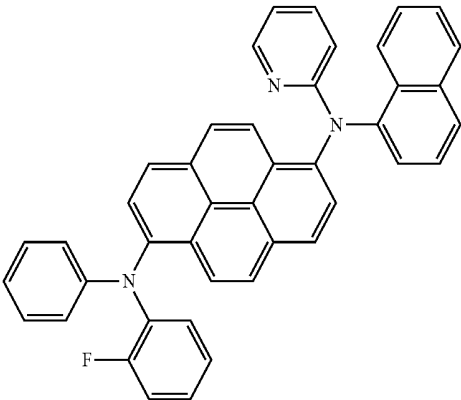
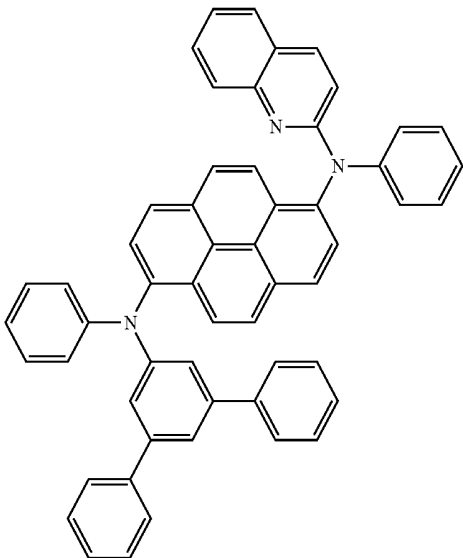
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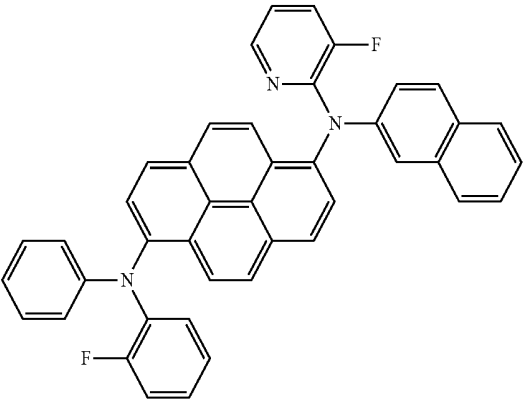
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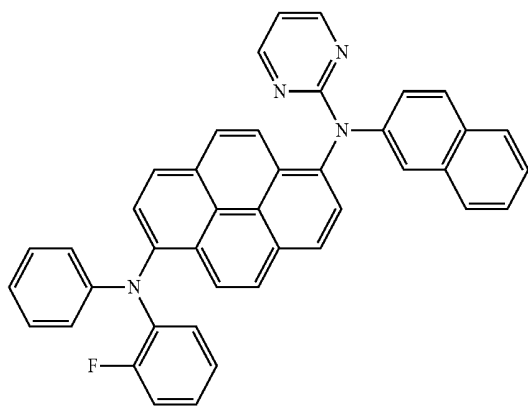


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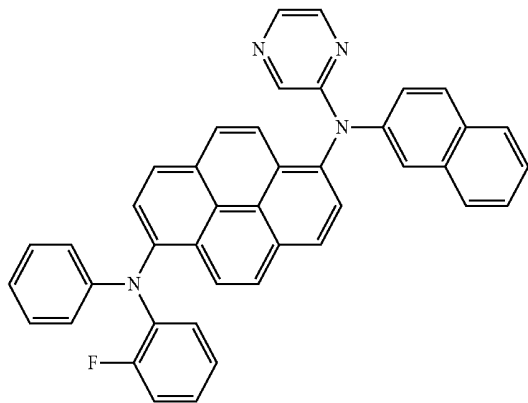


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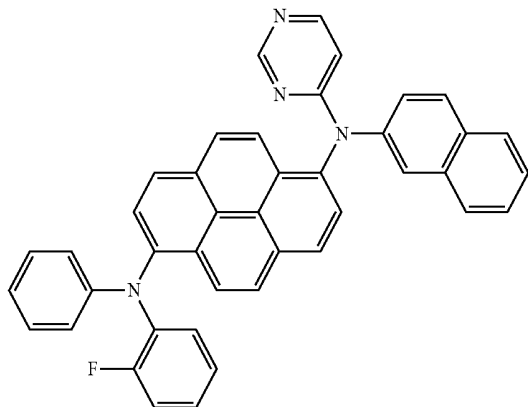
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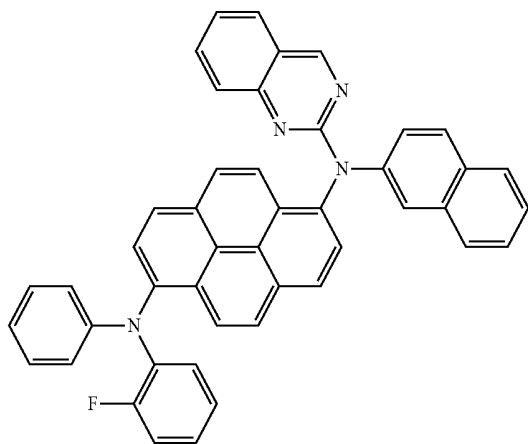
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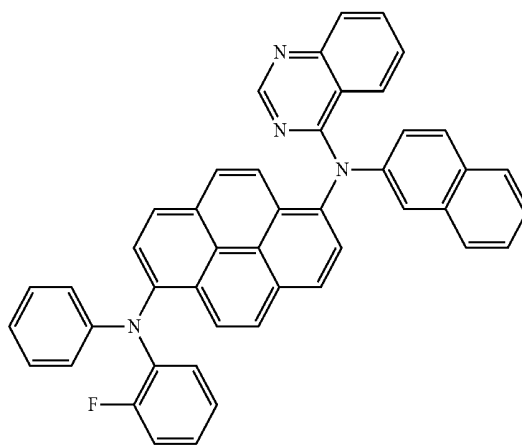


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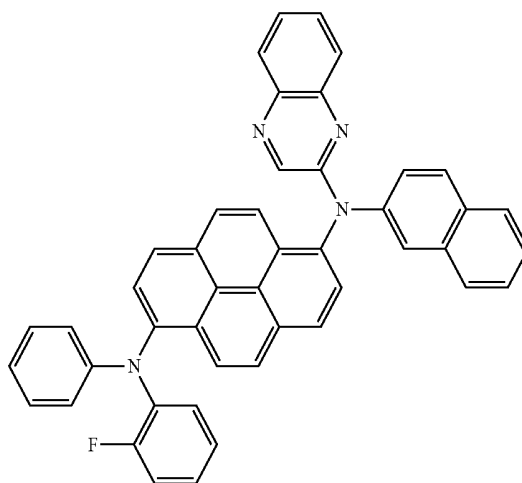


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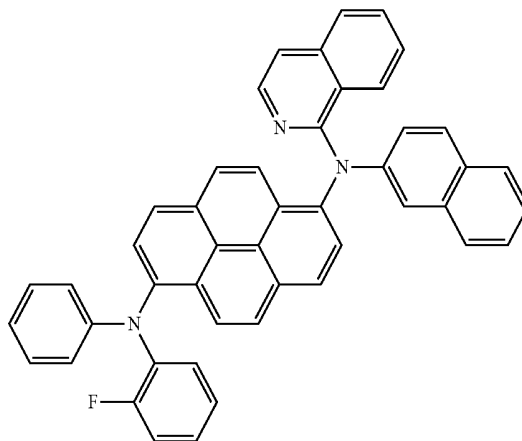
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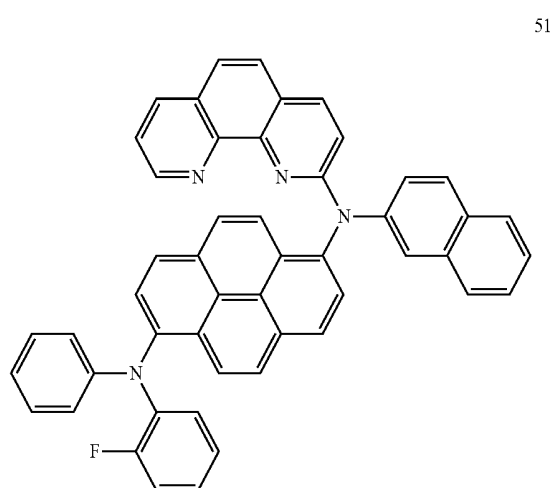
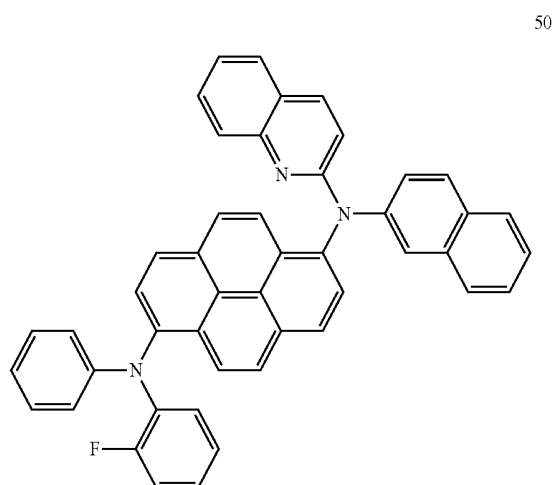
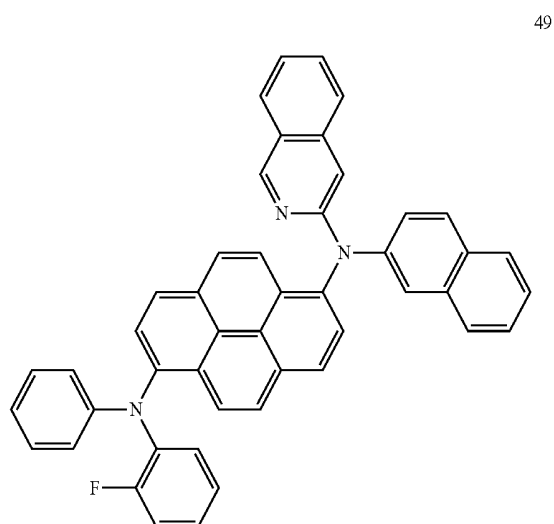
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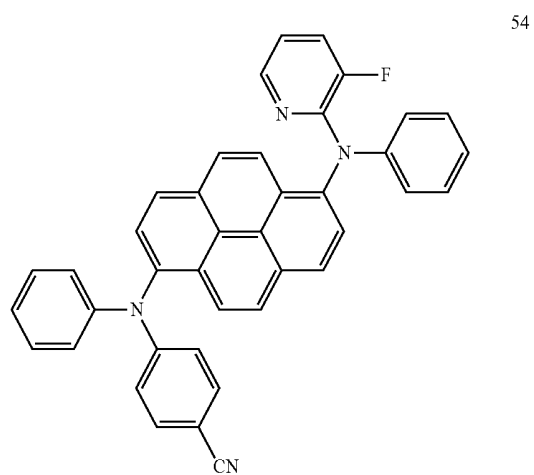
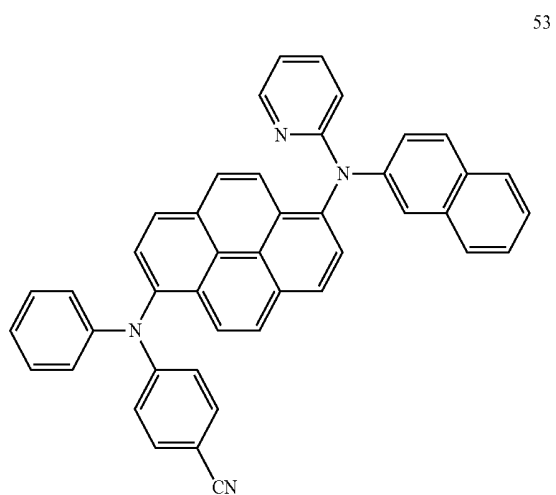
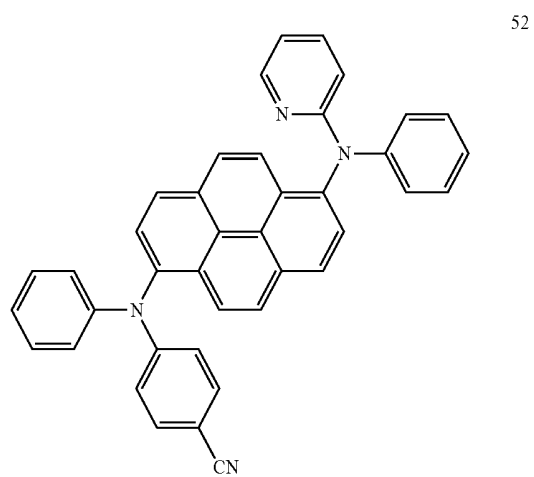
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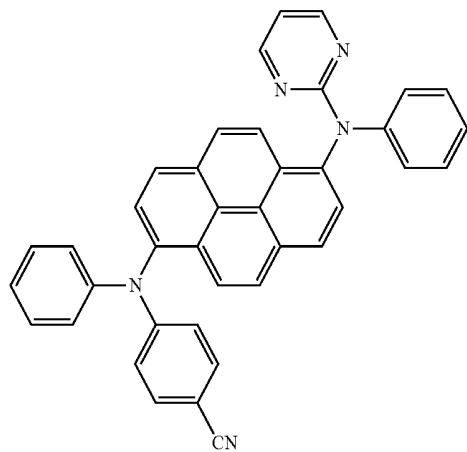
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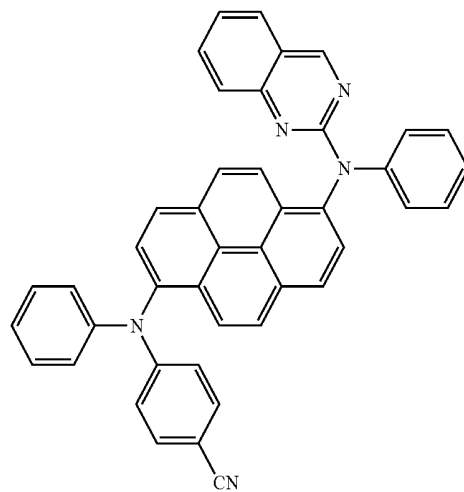


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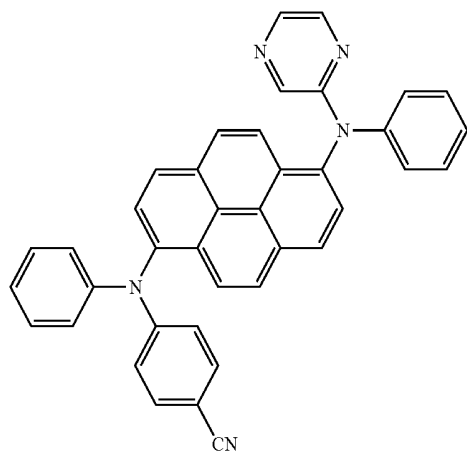


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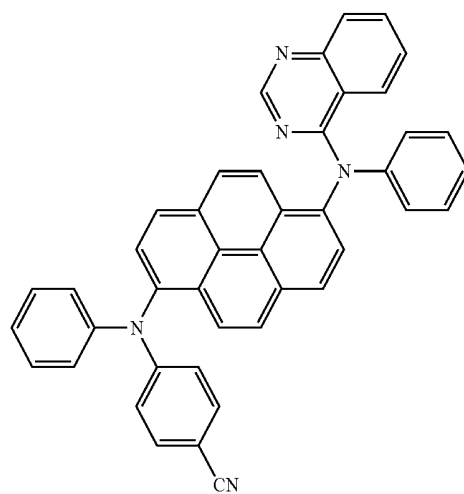
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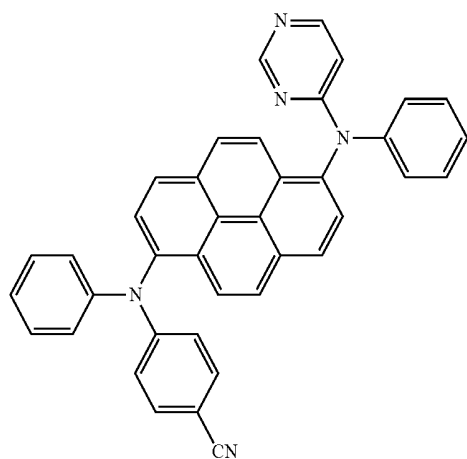
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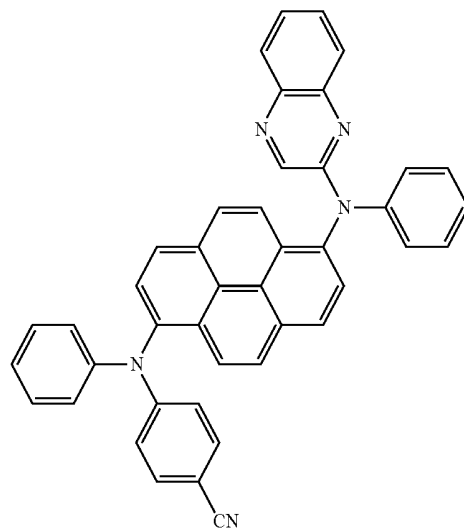
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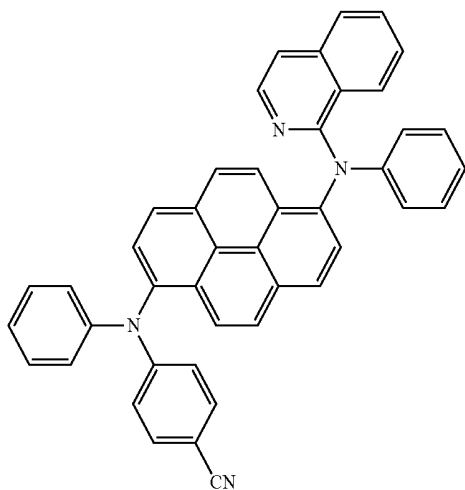


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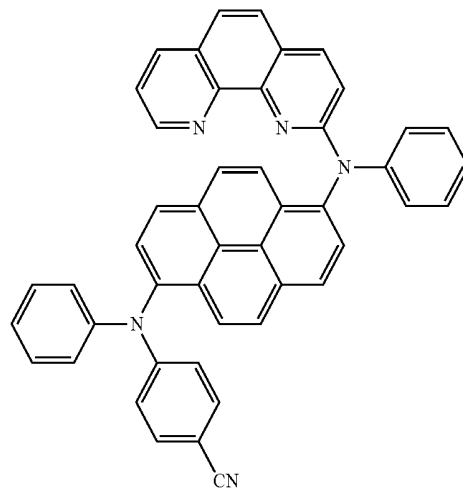
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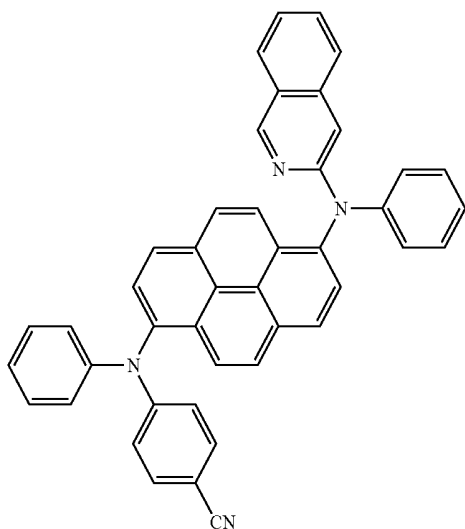


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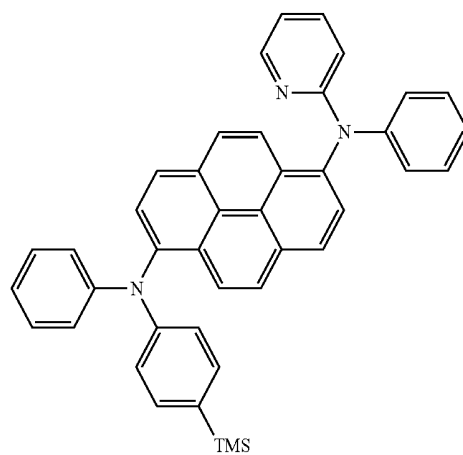
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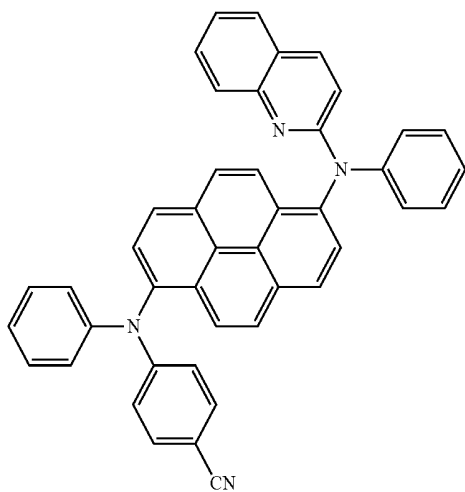
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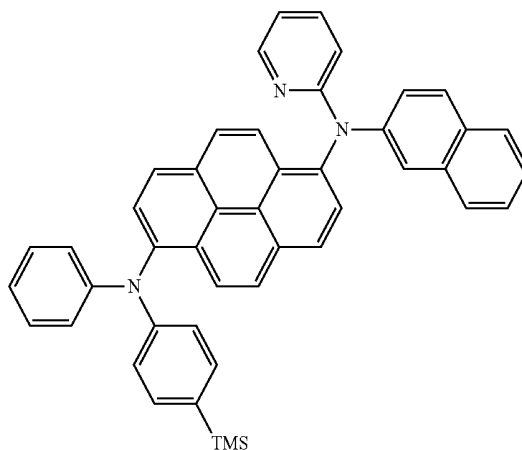
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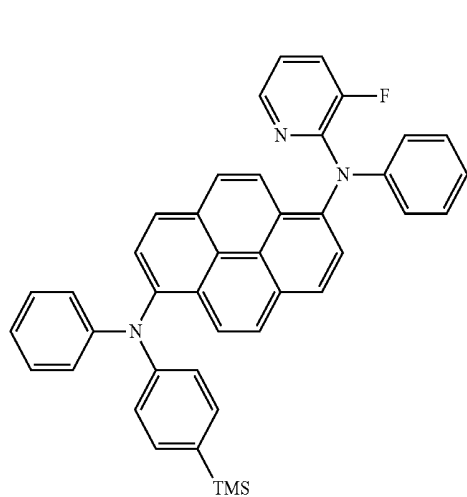


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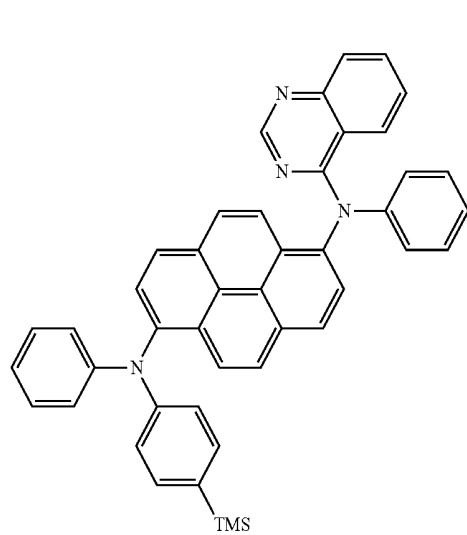
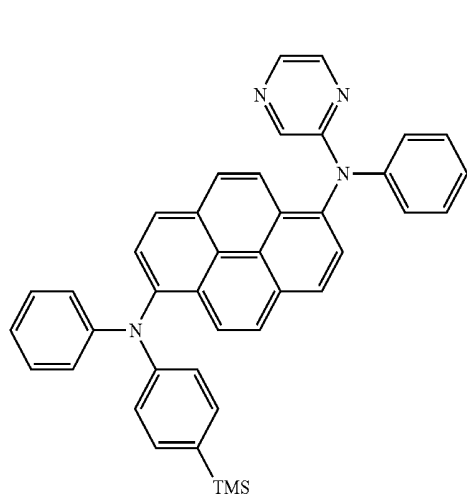
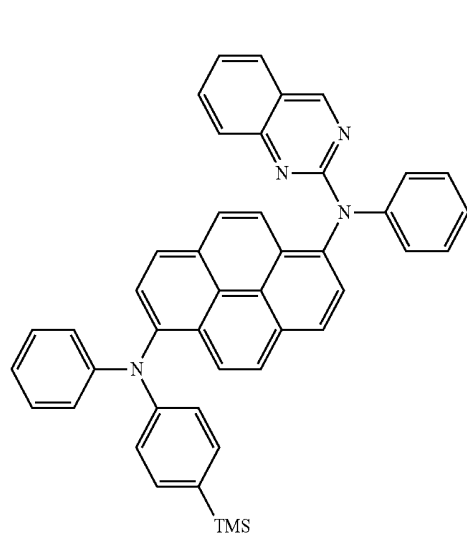
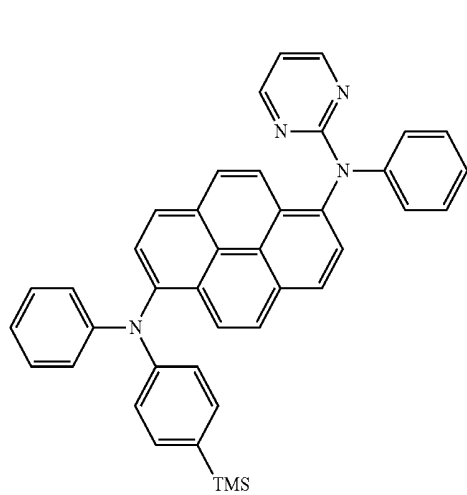
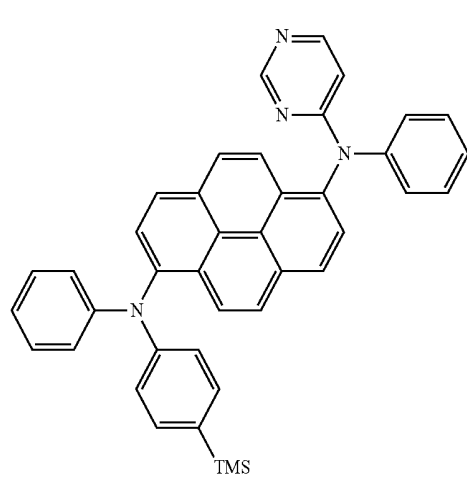


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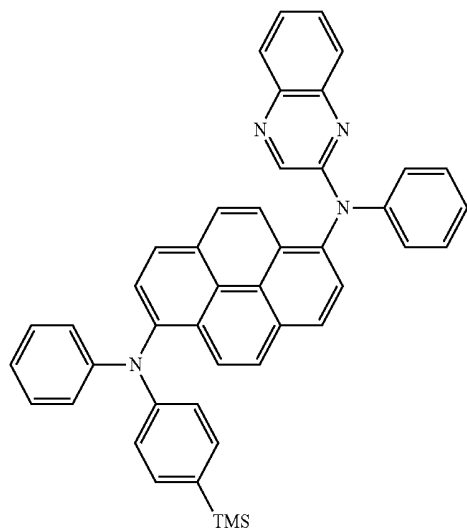


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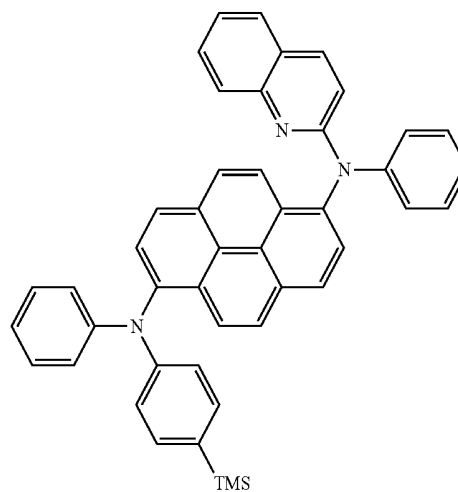
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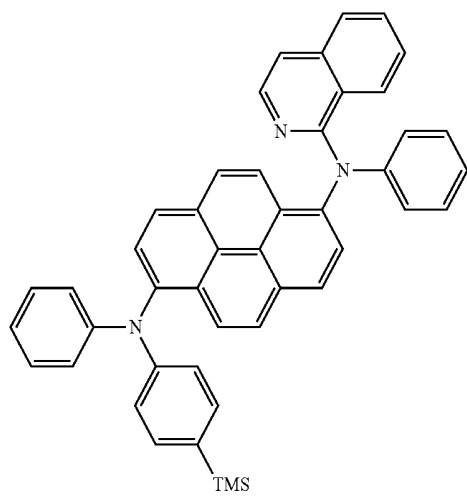


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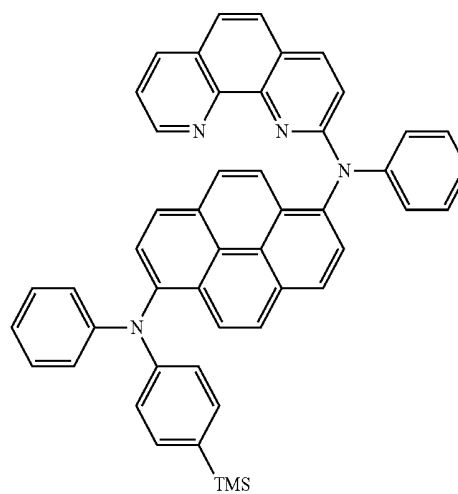
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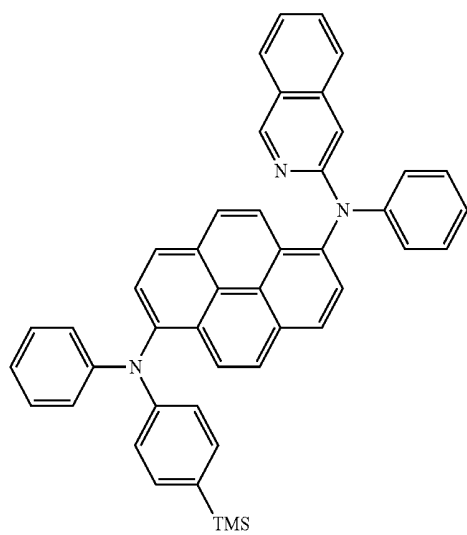
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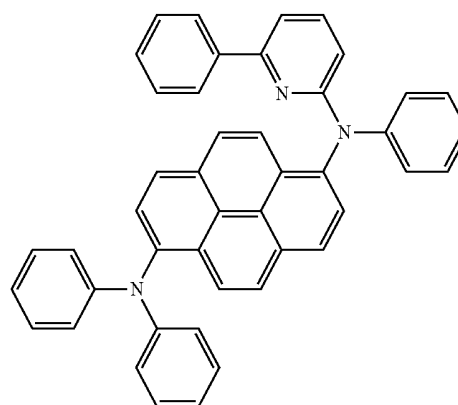
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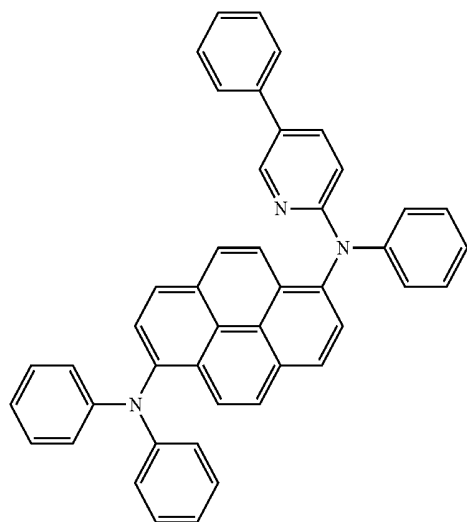
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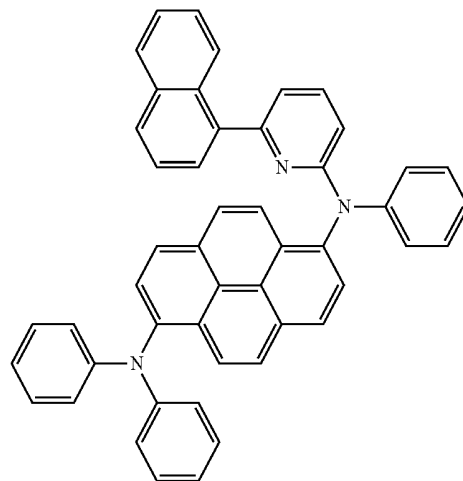


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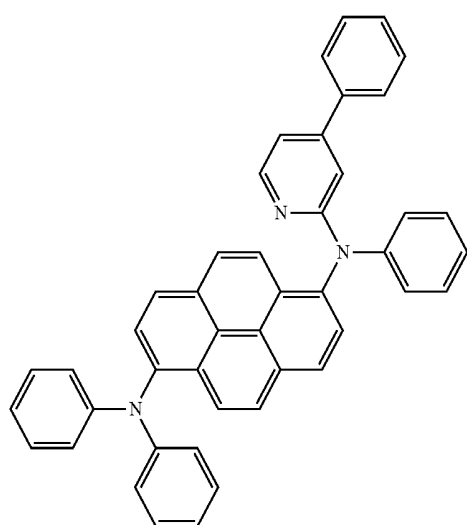


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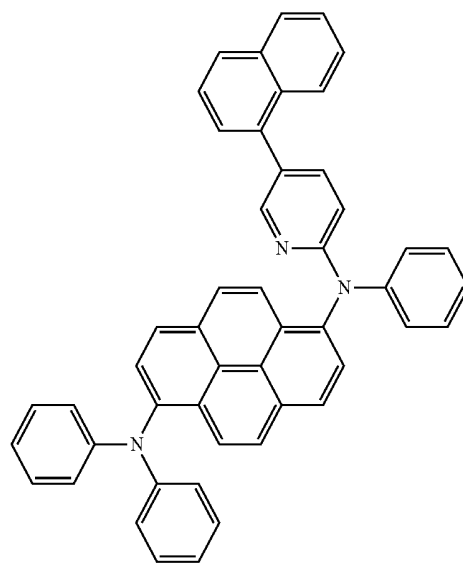
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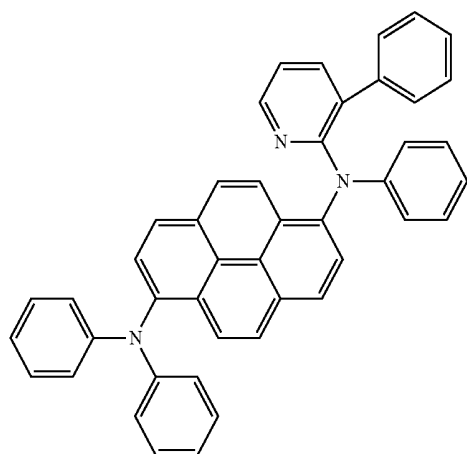
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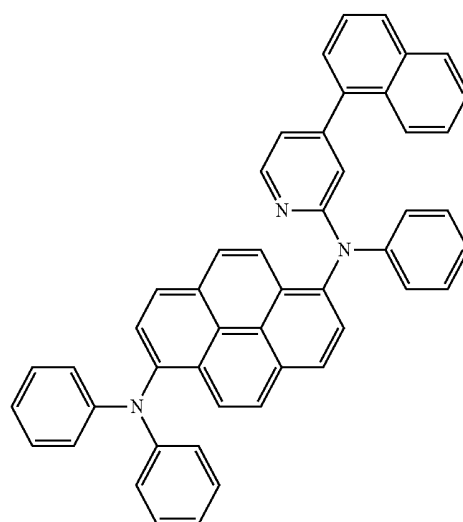
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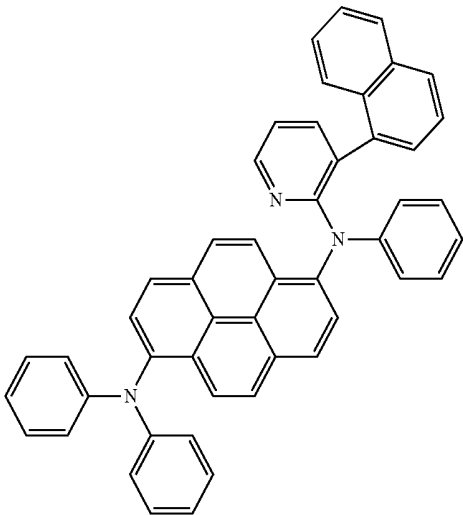


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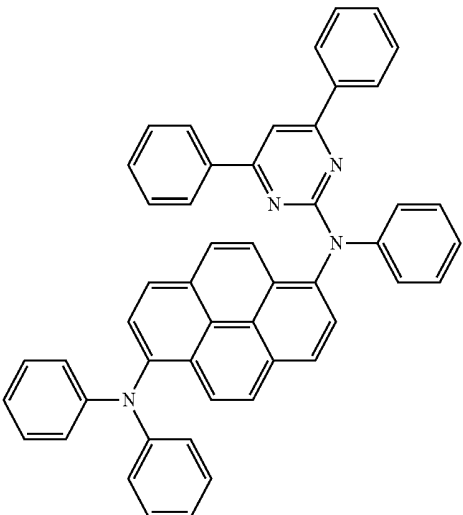
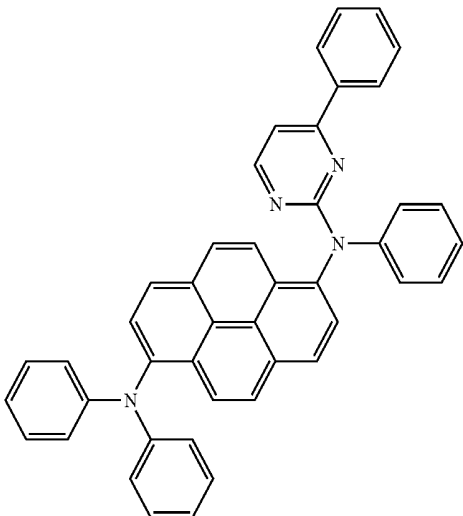
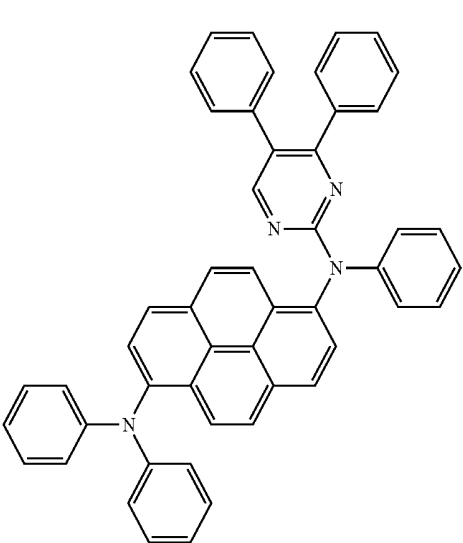
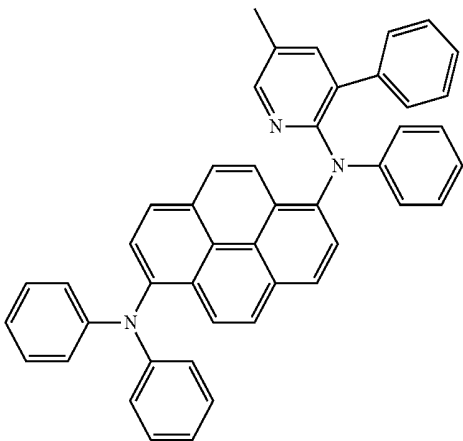
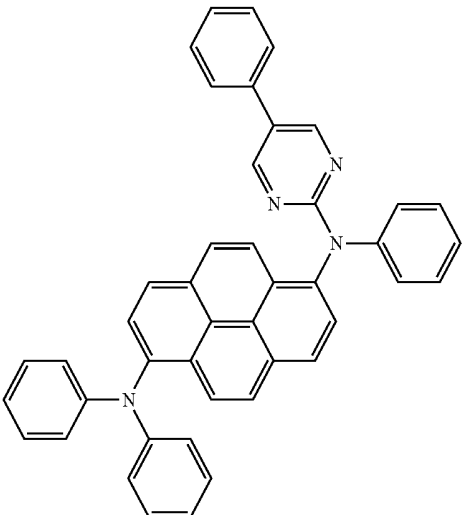


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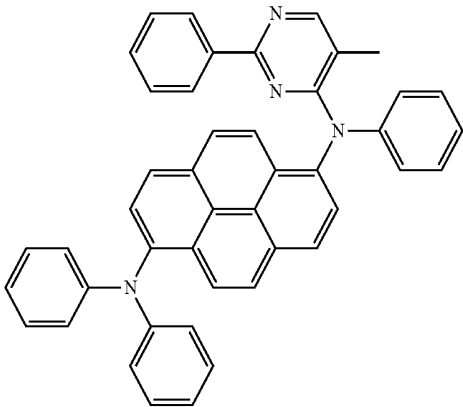
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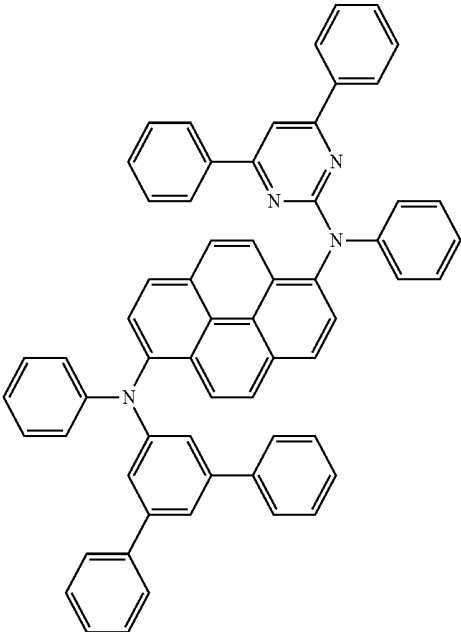


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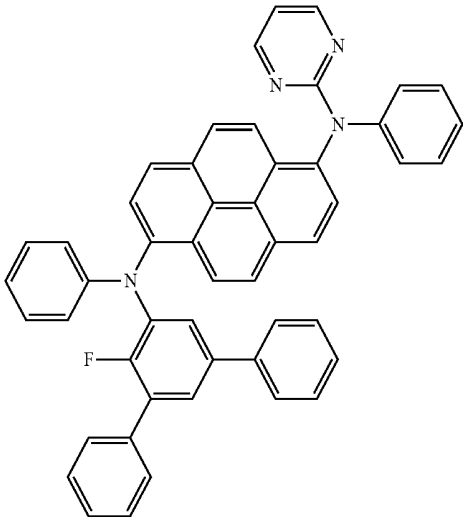


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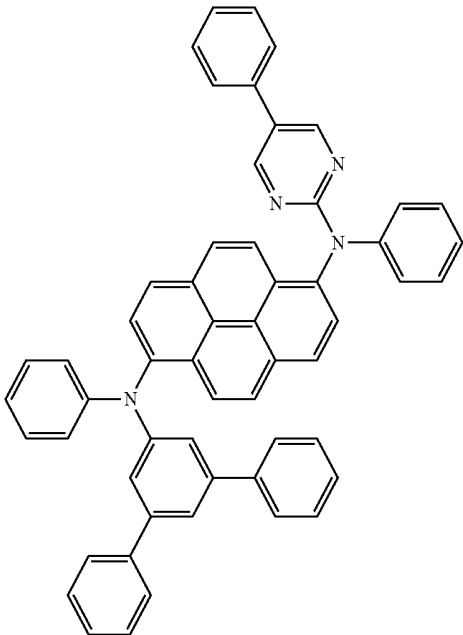
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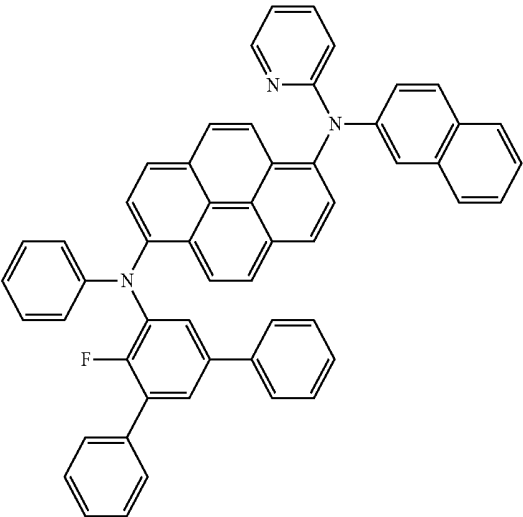
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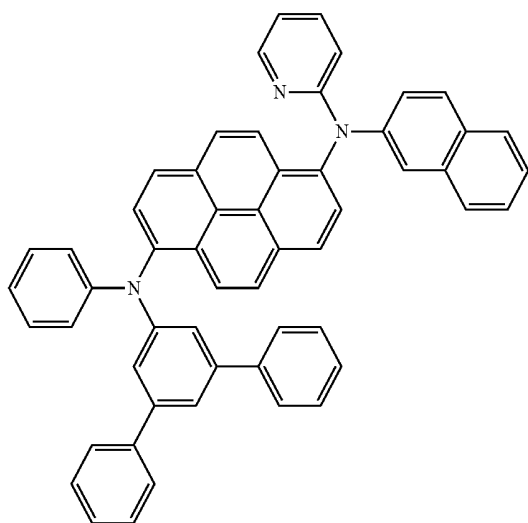
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[0092] The pyrene-based compound represented by Formula 1 may have, as a substituent of two amino groups, A, Ar₁, Ar₂, and Ar₃, which are described above. In this regard, A may be a heteroaromatic group including at least one N, and Ar₁, Ar₂ and Ar₃ may each be a substituted or unsubstituted phenyl group or a substituted or unsubstituted naphthyl group. Accordingly, the pyrene-based compound may be a heteroaromatic group in which only one (e.g., A) of four substituents (e.g., A, Ar₁, Ar₂ and Ar₃) of two amino groups has a hetero atom (e.g., nitrogen) as a ring element. Accordingly, the pyrene-based compound may emit blue light with excellent color purity characteristics. For example, an organic light-emitting diode including the pyrene-based compound represented by Formula 1 may provide blue light with high color purity of which a y coordinate is 0.11 or less.

[0093] Although not limited to a particular theory, in the case of an organic light-emitting diode including, as a dopant of an emission layer, a compound (for example, Compound B which will be describe below) that has the same structure as the pyrene-based compound represented by Formula 1 except that two or more of four substituents of two amino groups are heteroaromatic groups, energy transition from an anthracene-based host to the dopant in the emission layer may not sufficiently occur and thus, performance of the organic light-emitting diode during operation may be rapidly decreased.

[0094] Also, two amino groups of the pyrene-based compound represented by Formula 1 may be different from each other. Thus, the pyrene-based compound represented by Formula 1 may be asymmetric. Accordingly, a film including the pyrene-based compound represented by Formula 1 may have high amorphous properties, and thus, the film including the pyrene-based compound represented by Formula 1 may have high electric stability. Thus, an organic light-emitting diode using the pyrene-based compound represented by Formula 1 may provide high brightness at low current.

[0095] Accordingly, an organic light-emitting diode employing the pyrene-based compound of Formula 1 may provide excellent electric characteristics (e.g., a low driving voltage, a high current density, a long lifespan, or the like) and excellent color purity characteristics.

[0096] The pyrene-based compound represented by Formula 1 may be synthesized by using suitable organic synthesis methods. Such synthesis methods of a pyrene-based compound may be seen by referring to examples below.

[0097] One or more of the pyrene-based compound of Formula 1 may be used between a pair of electrodes of an organic light-emitting diode. For example, one or more of the pyrene-based compound may be used in an emission layer.

[0098] Accordingly, provided is an organic light-emitting diode according to an embodiment including: a first electrode; a second electrode facing the first electrode; and an organic layer interposed between the first electrode and the second electrode, wherein the organic layer includes one or more pyrene-based compounds described above.

[0099] The wording that“(an organic layer) includes at least one pyrene-based compound” used herein means that “(an organic layer) may include one pyrene-based compound represented by Formula 1 or two or more pyrene-based compounds represented by Formula 1 being different from each other”.

[0100] For example, the organic layer may include only Compound 3 as the pyrene-based compound. In this regard, Compound 3 may exist in an emission layer of the organic light-emitting diode. According to another embodiment, the organic layer may include Compound 3 and Compound 6 as the pyrene-based compound. In this regard, Compound 3 and Compound 6 may exist in an identical layer (for example, Compound 3 and Compound 6 may exist in an emission layer), or in different layers (for example, Compound 3 exists in an emission layer and Compound 6 exists in a hole transport layer.)

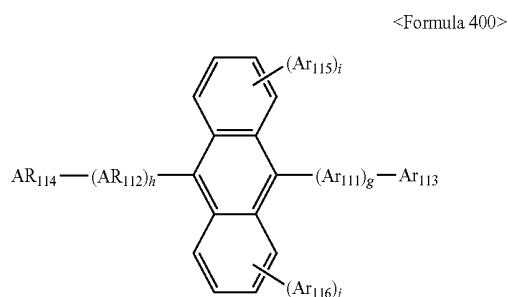
[0101] The organic layer may include at least one layer selected from a hole injection layer, a hole transport layer, a functional layer having a hole injection function and a hole transport function (hereinafter referred to as “H-functional layer”), a buffer layer, an electron blocking layer, an emission layer, a hole blocking layer, an electron transport layer, an electron injection layer, and/or a functional layer having an electron transport function and an electron injection function (hereinafter referred to as “E-functional layer”).

[0102] The term “organic layer” used herein refers to a single layer and/or a plurality of layers disposed between a first electrode and a second electrode of an organic light-emitting diode.

[0103] The organic layer may include an emission layer, and the emission layer may include one or more pyrene-based compounds described above.

[0104] The pyrene-based compound included in the emission layer may act as a dopant. For example, the pyrene-based compound may act as a fluorescent dopant. An emission layer including the pyrene-based compound may emit blue light. In this regard, the emission layer may further include a host.

[0105] The host may include at least one compound selected from an anthracene-based compound represented by Formula 400 below and an anthracene-based compound represented by Formula 401 below:

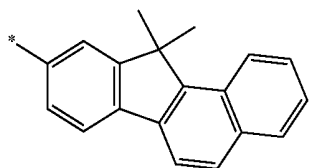


[0106] In Formula 400, Ar_{111} and Ar_{112} may each independently be a substituted or unsubstituted C_6 - C_{60} arylene group; Ar_{113} to Ar_{116} may each independently be a substituted or unsubstituted C_1 - C_{10} alkyl group or a substituted or unsubstituted C_6 - C_{60} aryl group; and g , h , i , and j may each independently be an integer of 0 to 4.

[0107] For example, Ar_{111} and Ar_{112} in Formula 400 may each independently be selected from a phenylene group, a naphthylene group, a phenanthrenylene group, or a pyrenylene group; and a phenylene group, a naphthylene group, a phenanthrenylene group, a fluorenyl group, and a pyrenylene group, each substituted with at least one of a phenyl group, a naphthyl group, and an anthracenyl group.

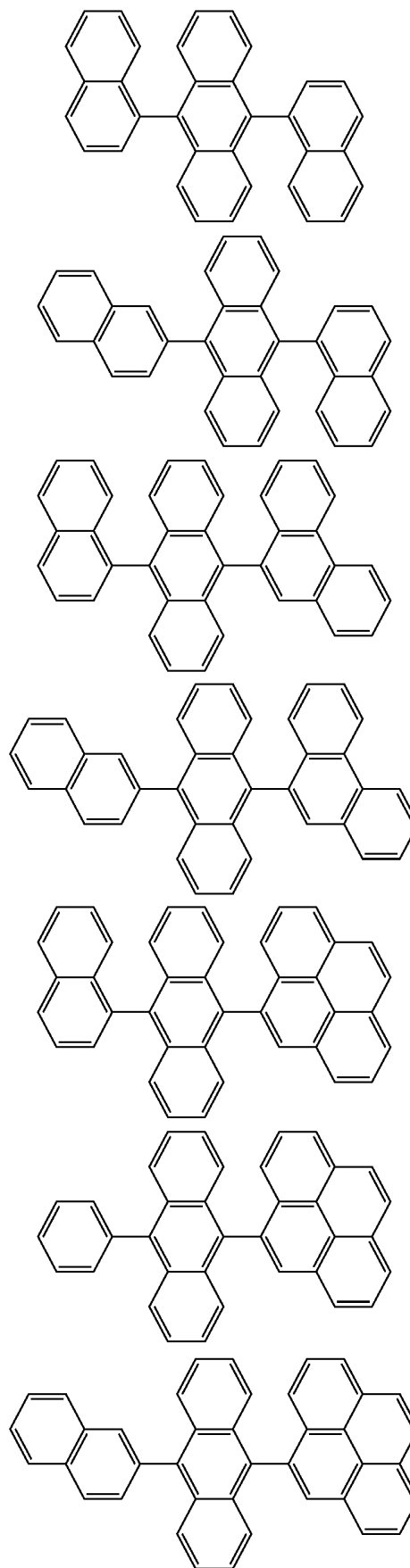
[0108] g , h , i , and j in Formula 400 may each independently be 0, 1, or 2.

[0109] Ar_{113} to Ar_{116} in Formula 400 may each independently be selected from a C_1 - C_{10} alkyl group, each substituted with at least one of a phenyl group, a naphthyl group, and an anthracenyl group; a phenyl group, a naphthyl group, an anthracenyl group, a pyrenyl group, a phenanthrenyl group and a fluorenyl group; and a phenyl group, a naphthyl group, an anthracenyl group, a pyrenyl group, a phenanthrenyl group, and a fluorenyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine, a hydrazone, a carboxyl group or a salt thereof, a sulfonic 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 anthracenyl group, a pyrenyl group,

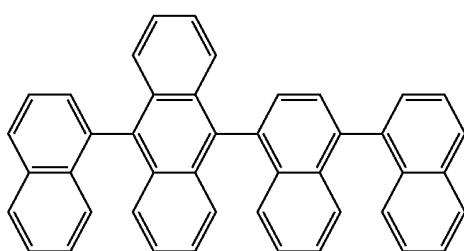
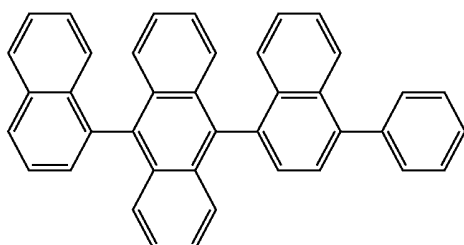
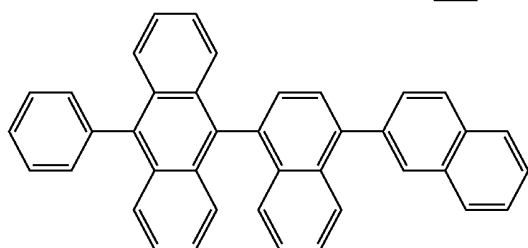
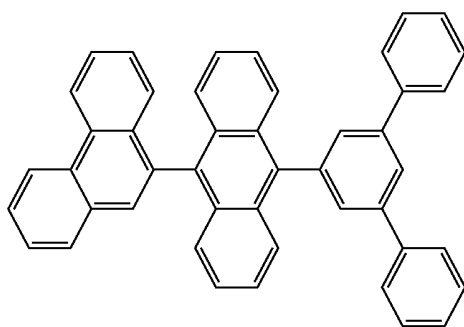
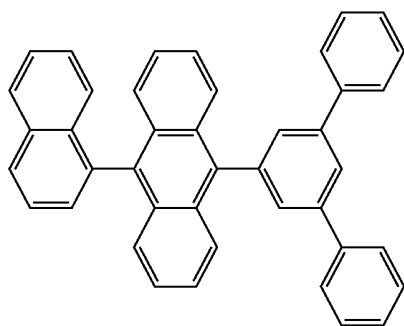
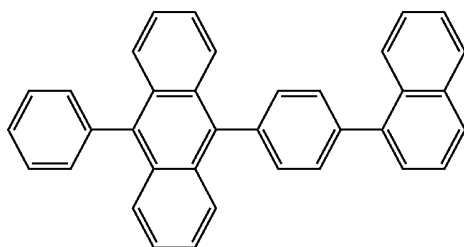


[0110] a phenanthrenyl group, a fluorenyl group, and

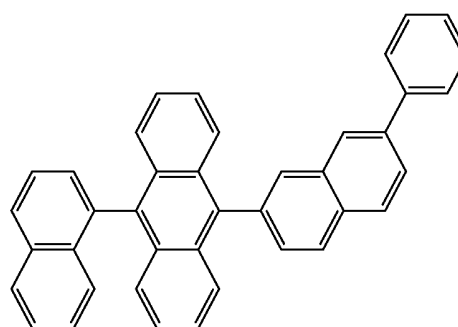
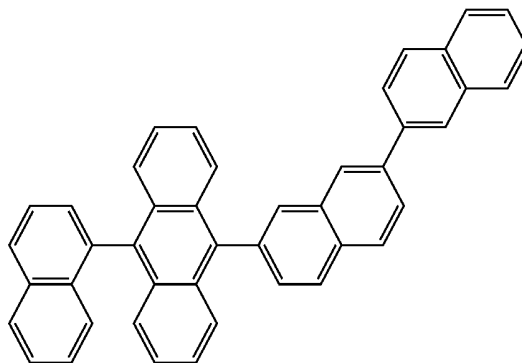
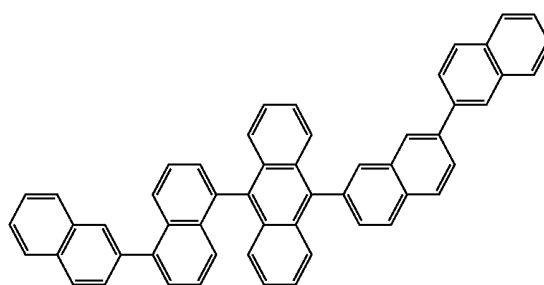
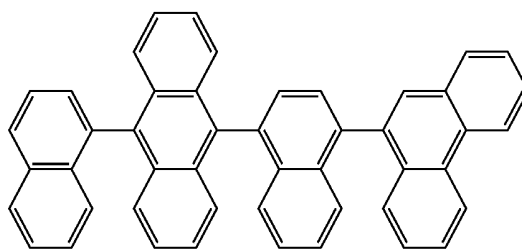
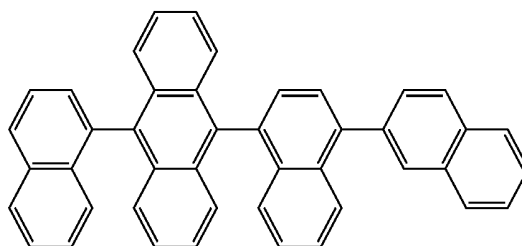
[0111] For example, the anthracene-based compound represented by Formula 400 may be one of the following compounds:



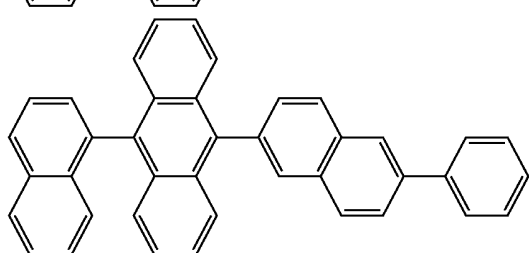
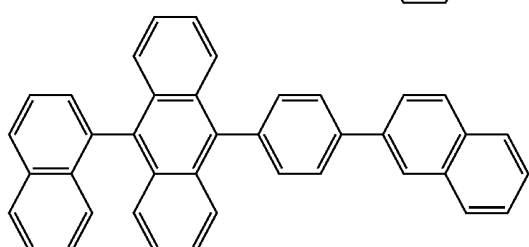
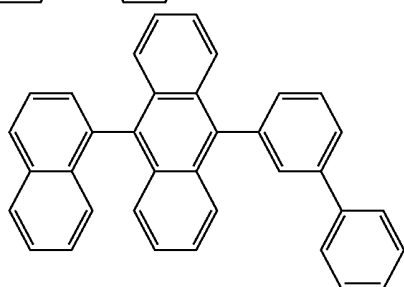
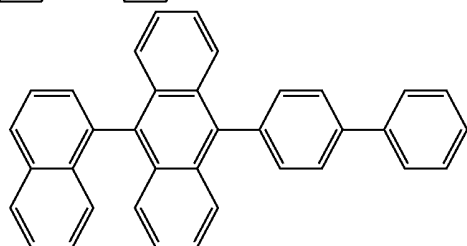
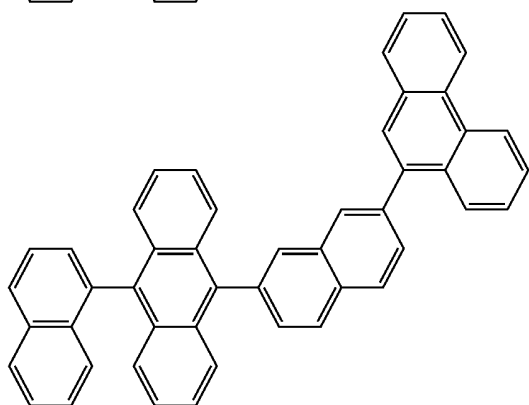
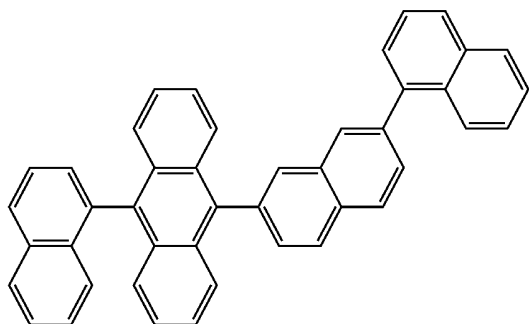
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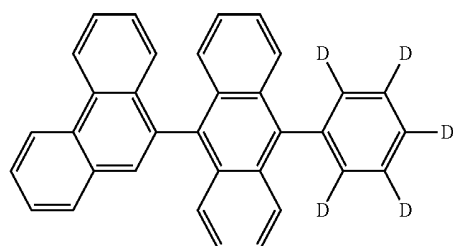
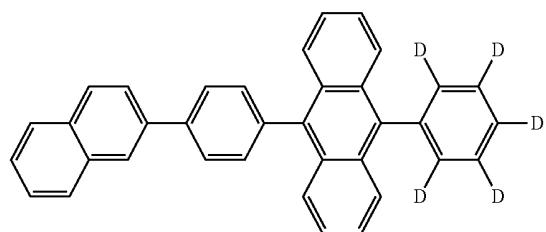
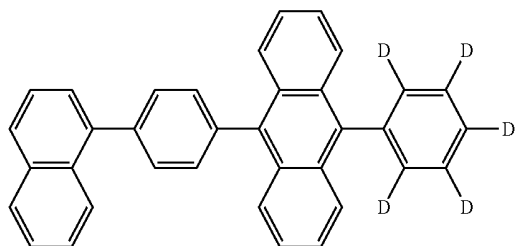
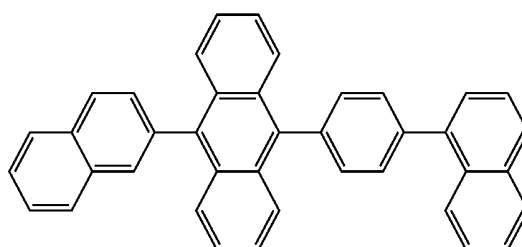
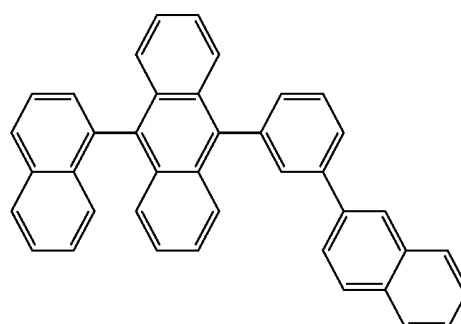
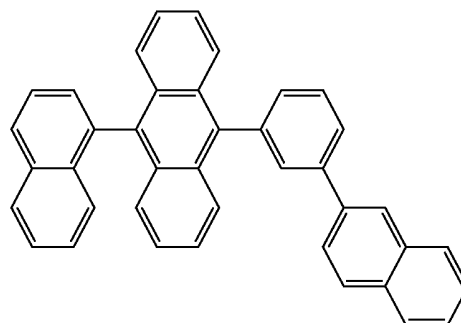
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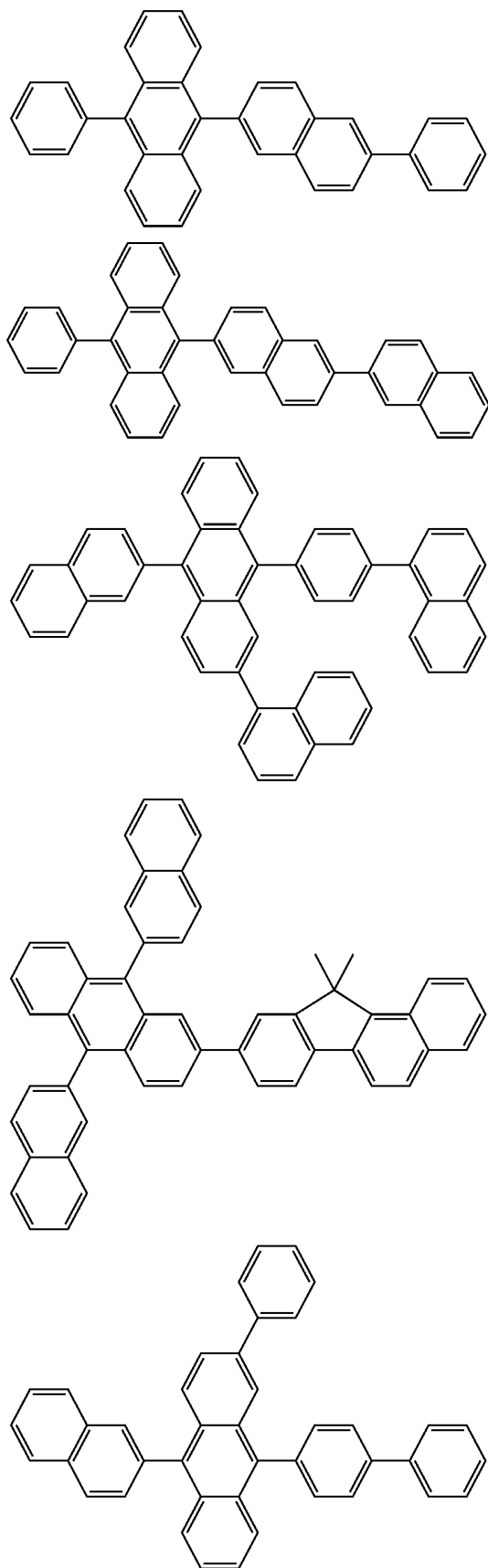
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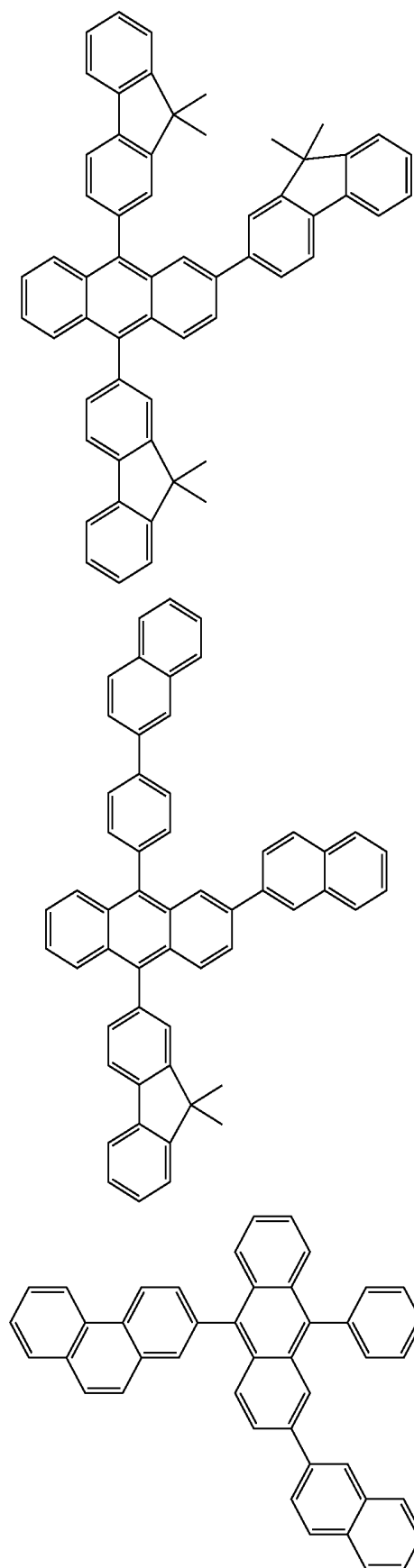
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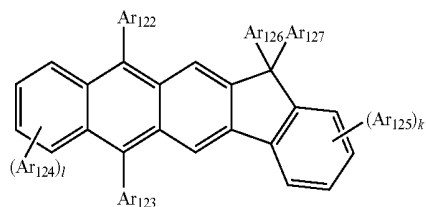


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<Formula 401>

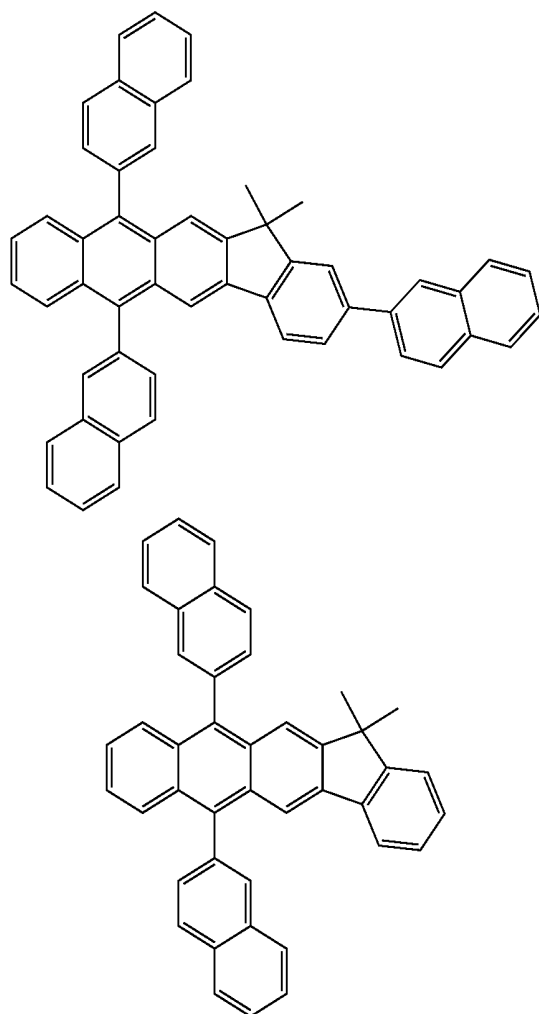


[0112] Ar_{122} to Ar_{125} in Formula 401 may be understood by referring to the description presented in connection to Ar_{113} in Formula 400.

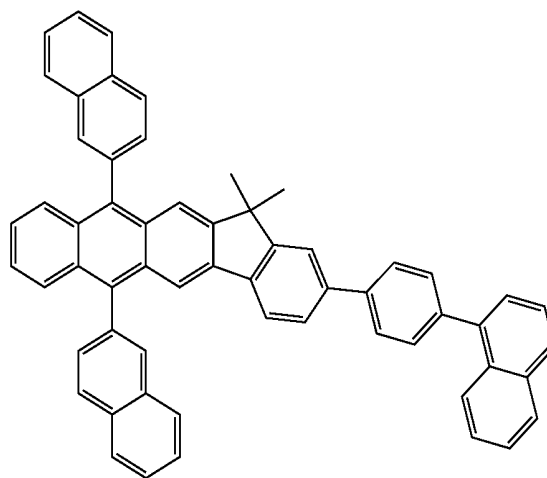
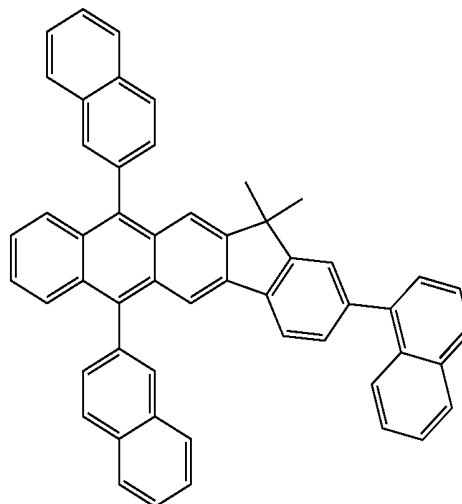
[0113] Ar_{126} and Ar_{127} in Formula 401 may each independently be a C_1 - C_{10} alkyl group (e.g., a methyl group, an ethyl group, or a propyl group).

[0114] k and l in Formula 401 may each independently be an integer of 0 to 4. For example, k and l may each independently be 0, 1, or 2.

[0115] For example, the anthracene-based compound represented by Formula 401 may be one of the following compounds.



-continued



[0116] An organic light-emitting diode employing the pyrene-based compound of Formula 1 may emit blue light that satisfies sRGB standard. Accordingly, the organic light-emitting diode may be used in a large full-color display device (e.g., an organic light-emitting display (OLED) TV, or the like.).

[0117] FIG. 1 illustrates a schematic sectional view of an organic light-emitting diode 10 according to an embodiment. Hereinafter, the structure of an organic light-emitting diode according to an embodiment and a method of manufacturing a pyrene-based compound according to an embodiment will be described in connection with FIG. 1.

[0118] A substrate 11, which may be a suitable substrate that is used for OLEDs, may be e.g., a glass substrate or a transparent plastic substrate, each with excellent mechanical strength, thermal stability, transparency, surface smoothness, ease of handling, and water resistance.

[0119] A first electrode 13 may be formed by depositing or sputtering a material for a first electrode on the substrate 11. When the first electrode 13 is an anode, the material for the first electrode may be selected from materials with a high work function to facilitate hole injection. The first electrode 13 may be a reflective electrode, a semi-transmissive electrode, or a transmissive electrode. For example, when the organic light-emitting diode is used for a large full-color display, the second electrode 13 of the organic light-emitting diode may be a reflective electrode or a semi-transmissive electrode. The material for the first electrode may be a transparent material with high conductivity, and examples of such a material may include indium tin oxide (ITO), indium zinc oxide (IZO), tin oxide (SnO₂), and zinc oxide (ZnO). When magnesium (Mg), aluminum (Al), aluminum-lithium (Al—Li), calcium (Ca), magnesium-indium (Mg—In), magnesium-silver (Mg—Ag), or the like is used, the first electrode 13 may be used as a reflective electrode.

[0120] 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.

[0121] An organic layer may be disposed of the first electrode 13.

[0122] The organic layer 15 may include a hole injection layer, a hole transport layer, a buffer layer, an emission layer, an electron transport layer, and an electron injection layer.

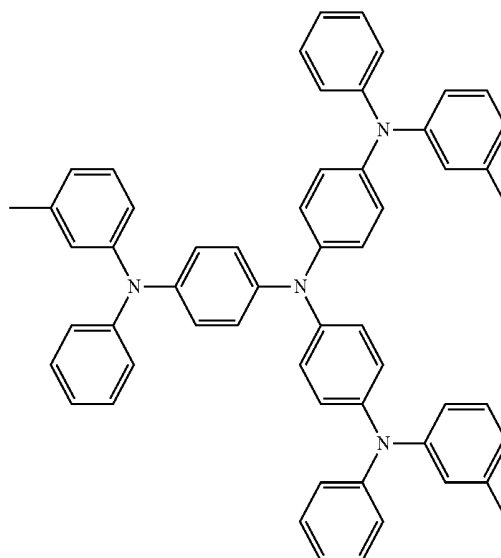
[0123] The hole injection layer (HIL) may be formed on the first electrode 13 by using various methods, e.g., vacuum deposition, spin coating, casting, Langmuir-Blodgett (LB) deposition.

[0124] 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 10⁻⁸ torr to about 10⁻³ torr, and a deposition rate of about 0.01 to about 100 Å/sec.

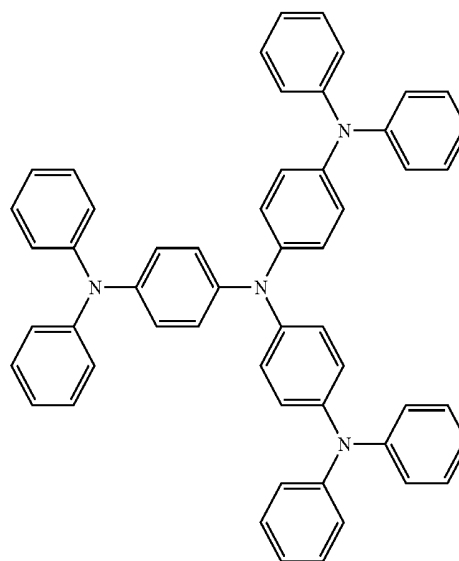
[0125] When a HIL is formed using spin coating, the coating conditions may vary according to the compound that is used to form the HIL, and the desired structure and thermal properties of the HIL to be formed. For example, the coating rate may be in the range of about 2,000 rpm to about 5,000 rpm, and 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.

[0126] For use as a hole injection material, a suitable hole injection material may be used. Examples of such a hole material may include 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), TDATA, 2-TNATA, polyaniline/dodecylbenzenesulfonic acid (pani/DBSA), poly(3,4-ethylenedioxy-

thiophene)/poly(4-styrenesulfonate) (PEDOT/PSS), polyaniline/camphor sulfonic acid (pani/CSA), or (polyaniline)/poly(4-styrenesulfonate) (PANI/PSS).

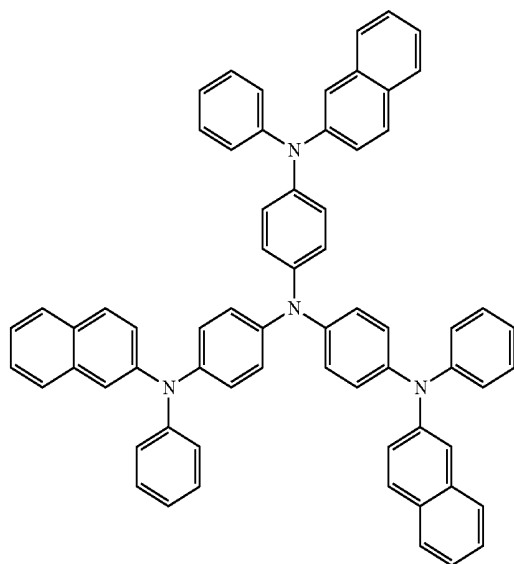


m-MTDATA



TDATA

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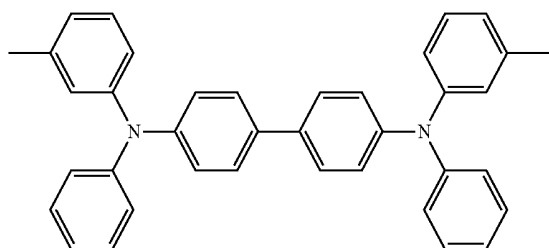


2-TNATA

[0127] A thickness of the HIL may be in a range of about 100 Å to about 10,000 Å, e.g., about 100 Å to about 1,000 Å. When the thickness of the HIL is within these ranges, the HIL may have satisfactory hole injection ability without a substantial increase in driving voltage.

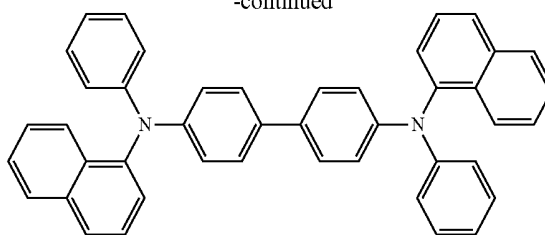
[0128] Then, a hole transportation layer (HTL) may be formed on the HIL by using vacuum deposition, spin coating, casting, Langmuir-Blodgett (LB) deposition, or the like. When a HTL is formed using vacuum deposition or spin coating, the deposition and coating conditions may be similar to those for the formation of the HIL, though the conditions for deposition and coating may vary according to the material that is used to form the HTL.

[0129] A material that is used to form the HTL may be a suitable hole transport material. Examples of the material that is used to form the HTL may include a carbazol derivative, such as N-phenylcarbazol or polyvinylcarbazol; N,N'-bis(3-methylphenyl)-N,N'-diphenyl-[1,1'-biphenyl]-4,4'-diamine (TPD); 4,4',4''-tris(N-carbazolyl)triphenylamine (TCTA), and alpha (α)-NPD.



TPD

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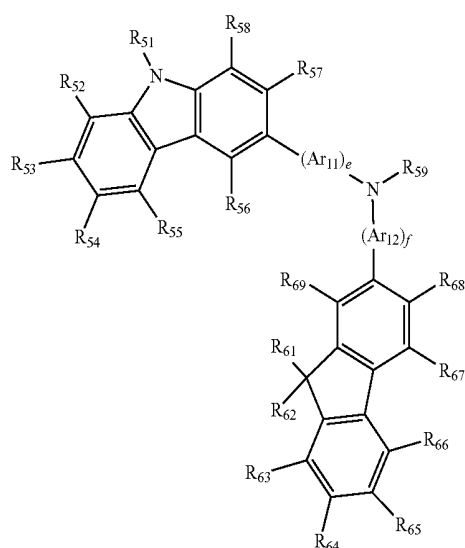
alpha (α)-NPD

[0130] A thickness of the HTL may be in a range of about 50 Å to about 20,000 Å, e.g., about 100 Å to about 1,500 Å. When the thickness of the HTL is within these ranges, the HTL may have satisfactory hole transporting ability without a substantial increase in driving voltage.

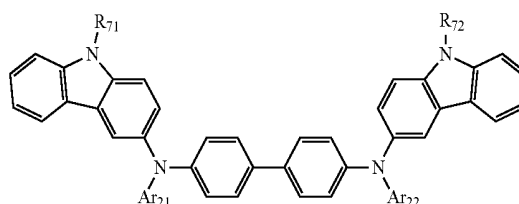
[0131] The H-functional layer (a functional layer having a hole injection ability and a hole transport ability) may include one or more materials selected from the materials for the HIL and the materials for the HTL. A thickness of the H-functional layer may be in a range of about 100 Å to about 10,000 Å, e.g., about 100 Å to about 1,000 Å. When the thickness of the H-functional layer is within these ranges, the H-functional layer may have satisfactory hole injection and transport characteristics without a substantial increase in driving voltage.

[0132] Also, at least one layer selected from the hole injection layer, the hole transport layer, and the H-functional layer may include at least one selected from a compound represented by Formula 300 below and a compound represented by Formula 350 below:

<Formula 300>



<Formula 350>



[0133] Ar_{11} and Ar_{12} in Formula 300 may each independently be a substituted or unsubstituted C_6-C_{60} arylene group.

[0134] For example, Ar_{11} and Ar_{12} may be each independently selected from:

[0135] a phenylene group, a pentalenylene group, an indenylene group, a naphthylene group, an azulenylene group, an acenaphthylene group, a fluorenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylene group, a fluorantenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylenylene group, a naphthacenylene group, a pycenylene group, a perylenylene group and a pentacenylene group; and

[0136] a phenylene group, a pentalenylene group, an indenylene group, a naphthylene group, an azulenylene group, an acenaphthylene group, a fluorenylene group, a phenalenylene group, a phenanthrenylene group, an anthracenylene group, a fluorantenylene group, a triphenylenylene group, a pyrenylene group, a chrysenylenylene group, a naphthacenylene group, a pycenylene group, a perylenylene group and a pentacenylene group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic 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 C_3-C_{10} cycloalkyl group, a C_3-C_{10} cycloalkenyl group, a C_3-C_{10} heterocycloalkyl group, a C_3-C_1 heterocycloalkenyl group, a C_6-C_{60} aryl group, a C_6-C_{60} aryloxy group, a C_6-C_{60} arylthio group, and a C_2-C_{60} heteroaryl group.

[0137] e and f in Formula 300 may each independently be an integer of 0 to 5, or 0, 1, or 2. In an implementation, e may be 1 and f may be 0.

[0138] R_{51} to R_{58} , R_{61} to R_{69} and R_{71} to R_{72} in Formulae 300 and 350 may each independently be selected from a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic 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_{60} aryl group, a substituted or unsubstituted C_6-C_{60} aryloxy group, and a substituted or unsubstituted C_6-C_{60} arylthio group.

[0139] For example, R_{51} to R_{58} , R_{61} to R_{69} , and R_{71} to R_{72} may each independently be selected from:

[0140] a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic 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, or a hexyl group), and 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);

[0141] a C_1-C_{10} alkyl group and a C_1-C_{10} alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof;

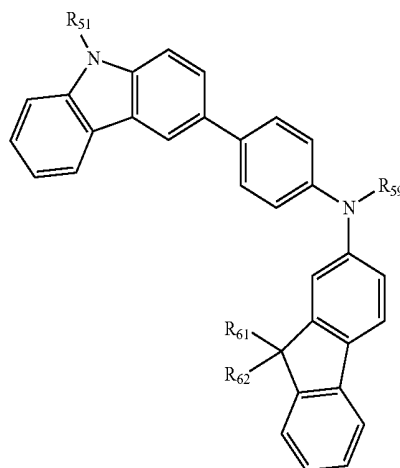
[0142] a phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, and a pyrenyl group; and

[0143] a phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, and a pyrenyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C_1-C_{10} alkyl group, and a C_1-C_{10} alkoxy group, and is not limited thereto.

[0144] R_{59} , Ar_{21} , and Ar_{22} in Formulae 300 and 350 may each independently be selected from a phenyl group, a naphthyl group, an anthracenyl group, biphenyl group, and a pyridyl group; and a phenyl group, a naphthyl group, an anthracenyl group, biphenyl group, and a pyridyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic 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.

[0145] In an implementation, the compound represented by Formula 300 may be represented by Formula 300A below.

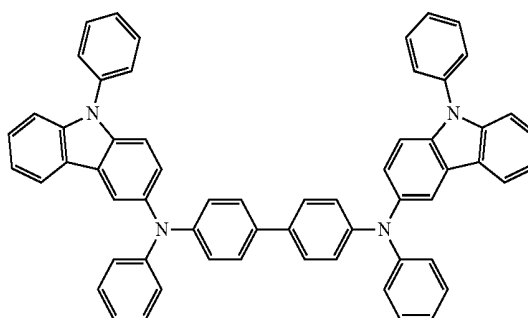
<Formula 300A>



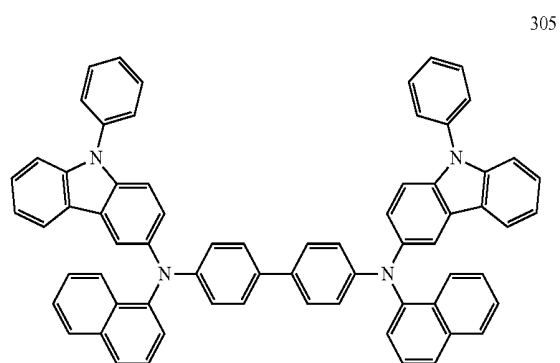
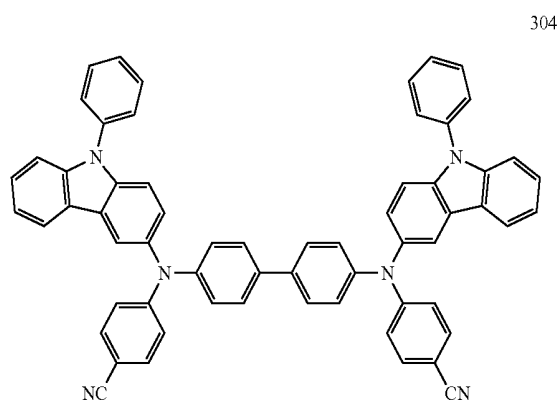
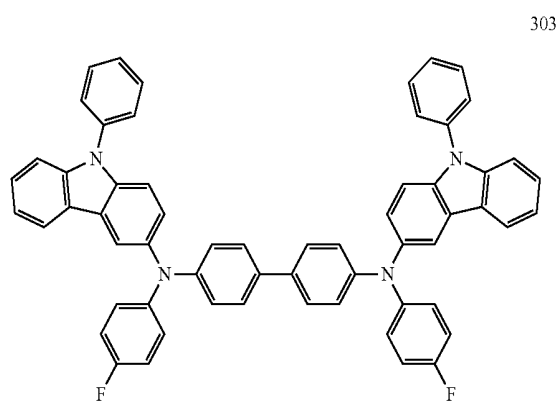
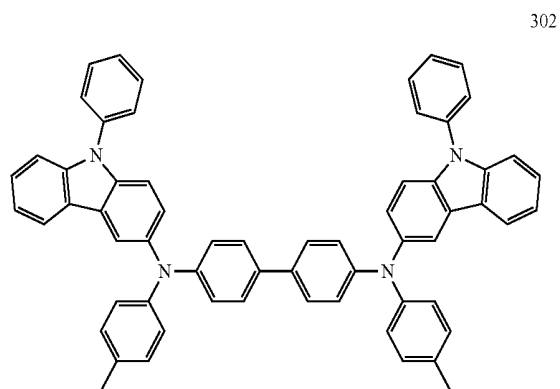
[0146] R_{51} , R_{61} , R_{62} and R_{59} in Formula 300A are already described in detail above.

[0147] For example, at least one layer selected from the HIL, the HTL, and the H-functional layer may include at least one selected from Compounds 301 to 320. In an implementation, these layers may instead or additionally include other compounds:

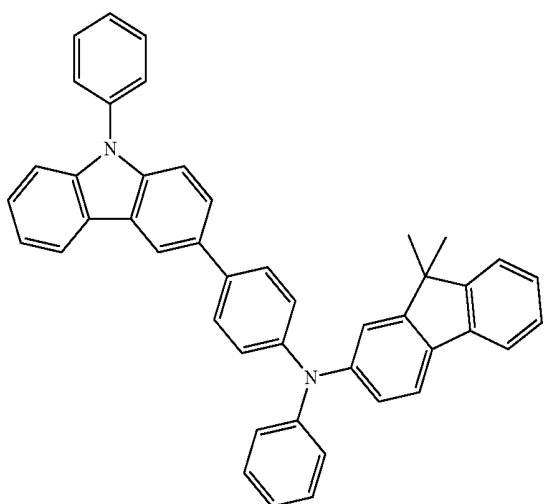
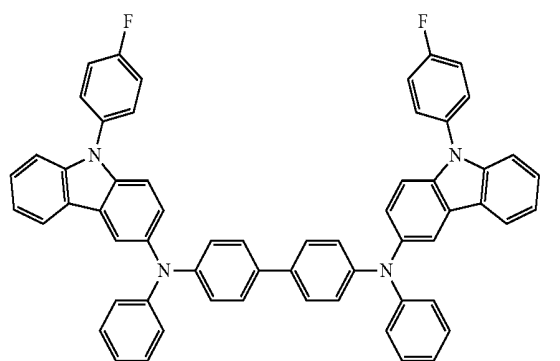
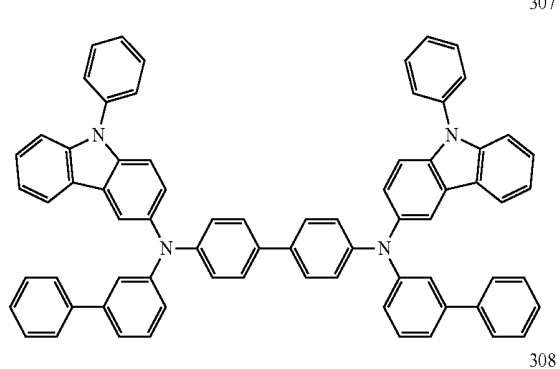
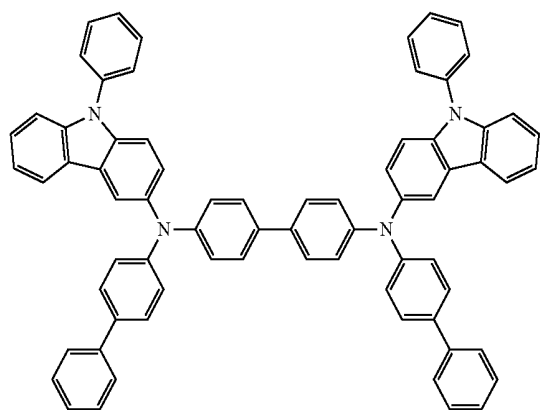
301



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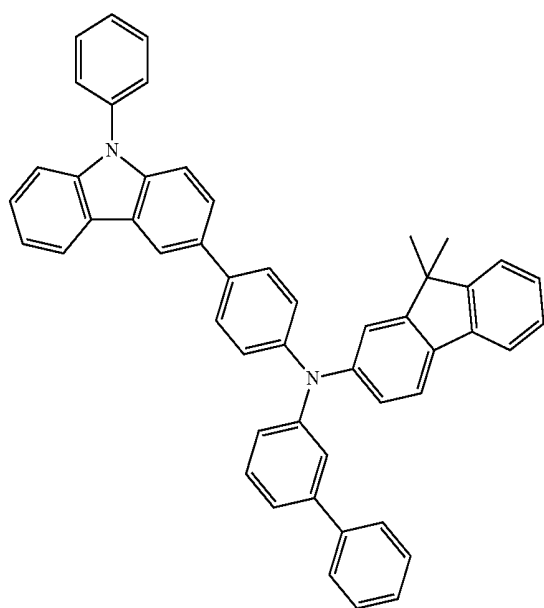
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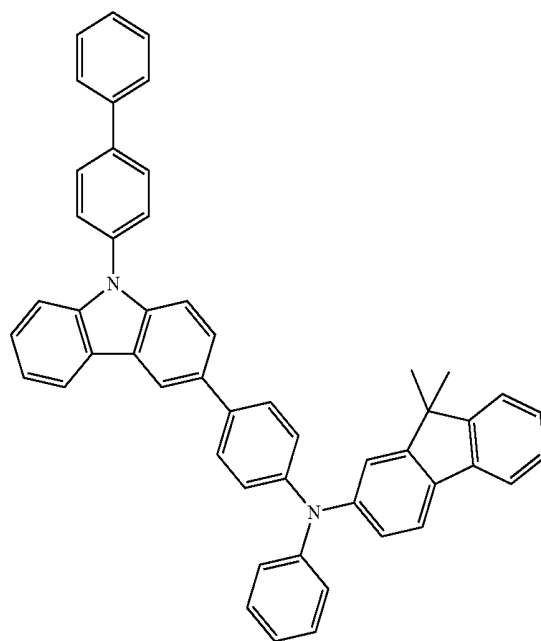
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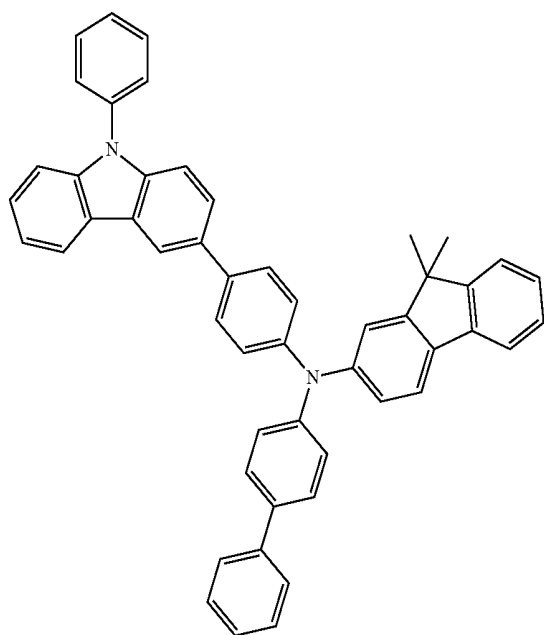
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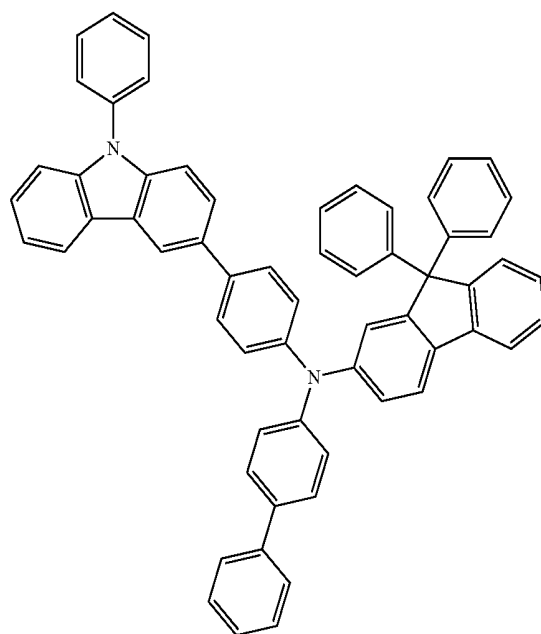
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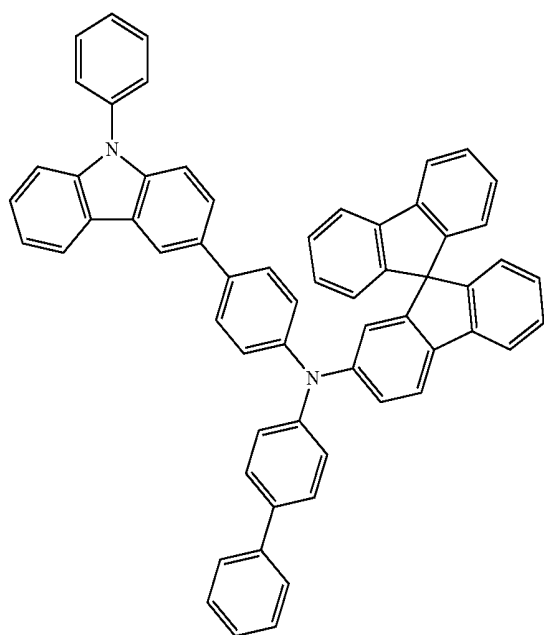
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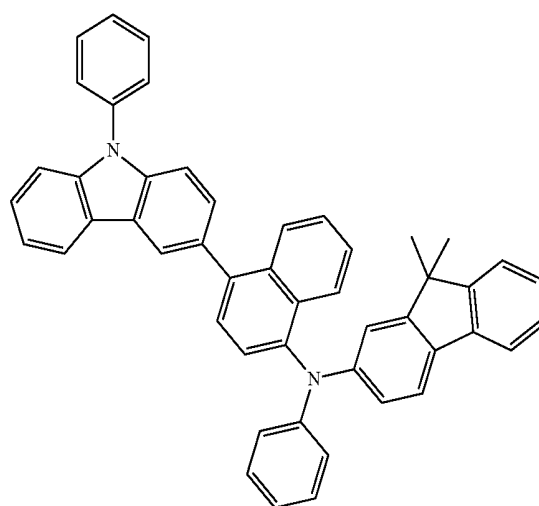
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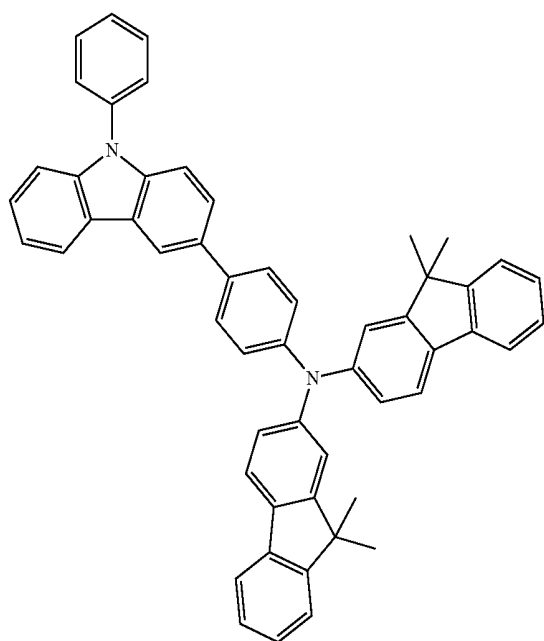
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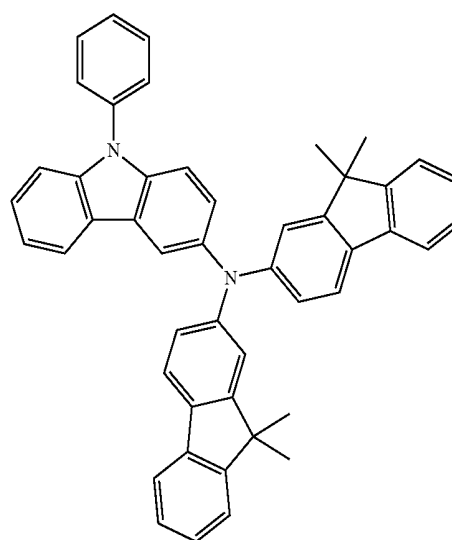
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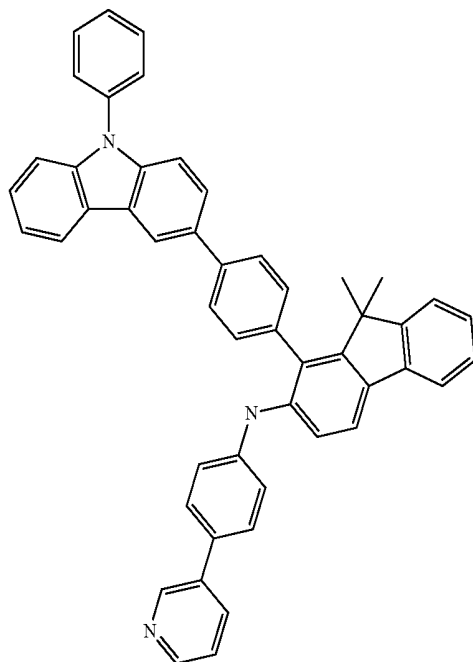
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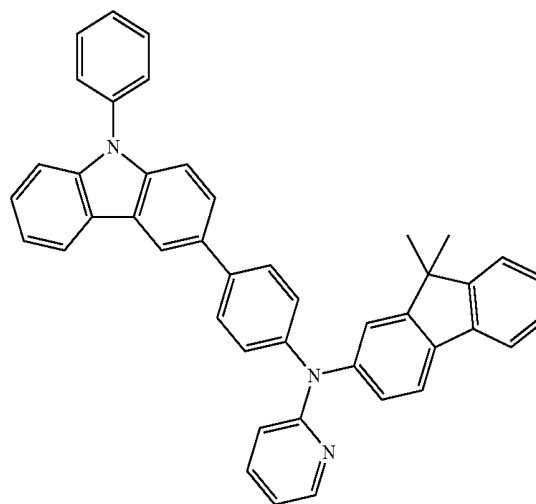


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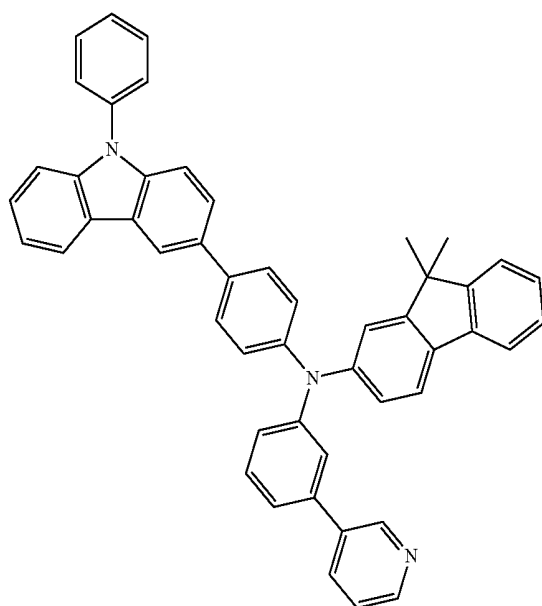
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320

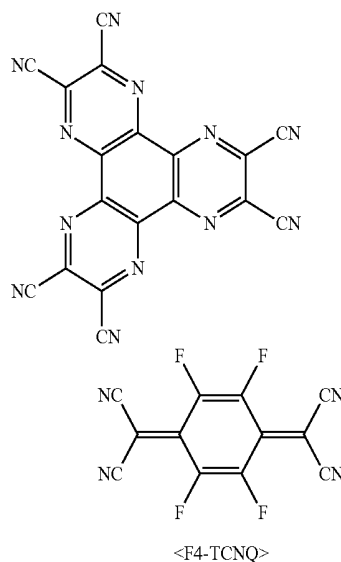


[0148] At least one layer selected from the hole injection layer, the hole transport layer, and the H-functional layer may further include, in addition to suitable hole injection materials, suitable hole transport materials, and/or suitable materials having a hole injection function and a hole transport function, a charge-generating material to help increase conductivity of the corresponding layer.

[0149] The charge-generating material may be, e.g., a p-dopant. The p-dopant may be one of quinine derivatives, metal oxides, and compounds with a cyano group, but is not limited thereto. Non-limiting examples of the p-dopant are a quinone derivative, such as tetracyanoquinonodimethane (TCNQ) or 2,3,5,6-tetrafluoro-tetracyano-1,4-benzoquinonodimethane (F4-TCNQ); a metal oxide, such as tungsten oxide or molybdenum oxide; and a cyano group-containing compound, such as Compound 200 below.



<Compound 200>



[0150] When the HIL, the HTL, or the H-functional layer further includes a charge-generating material, the charge-generating material may be homogeneously dispersed or non-homogeneously distributed in the HIL, the HTL, or the H-functional layer.

[0151] A buffer layer may be disposed between the EML and at least one of the HIL, the HTL, and the H-functional layer. The buffer layer may help compensate for an optical resonance distance of light according to a wavelength of the light emitted from an emission layer, and thus may help increase efficiency. The buffer layer may include suitable hole injecting materials or hole transporting materials. In an implementation, the buffer layer may include the same material as one of the materials included in the HIL, the HTL, and the H-functional layer that are disposed under the buffer layer.

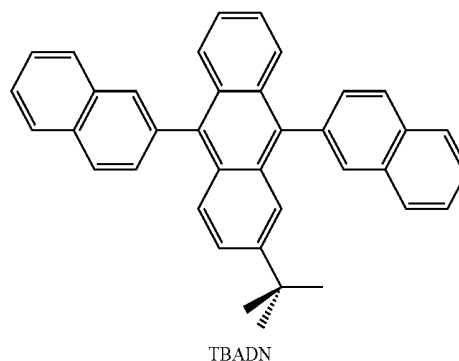
[0152] Then, an emission layer (EML) may be formed on the HIL, the HTL, or the H-functional layer by, e.g., vacuum deposition, spin coating, casting, LB deposition, or the like. When the EML is formed using vacuum deposition or spin coating, the deposition and coating conditions may be similar to those 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.

[0153] The EML may include one or more of the pyrene-based compounds described above.

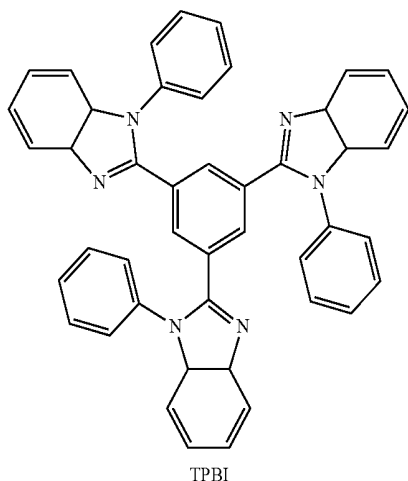
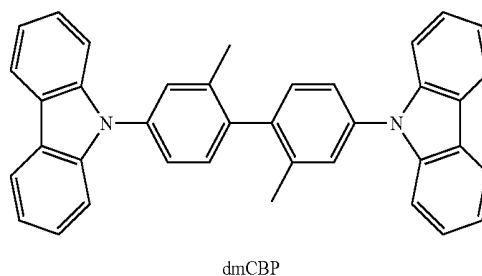
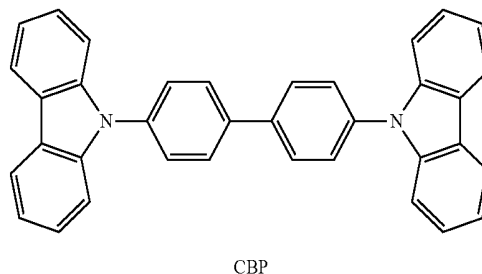
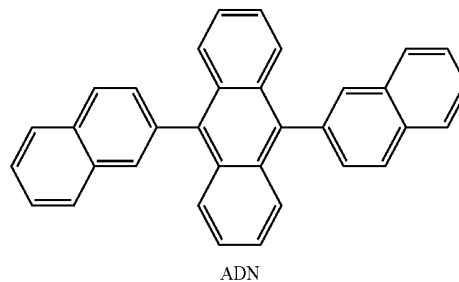
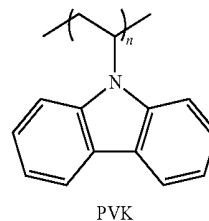
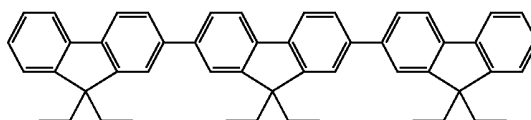
[0154] The pyrene-based compound included in the EML may act as a dopant (e.g., a blue fluorescent dopant). For example, the EML may further include, in addition to the pyrene-based compound, a host.

[0155] As the host, Alq₃, 4,4'-N,N'-dicarbazol-biphenyl (CBP), poly(n-vinylcarbazole) (PVK), 9,10-di(naphthalen-2-yl)anthracene (ADN), TCTA, 1,3,5-tris(N-phenylbenzimidazole-2-yl)benzene (TPBI), 3-tert-butyl-9,10-di(naphth-2-yl) anthracene (TBADN), E3, distyryl arylene (DSA), dmCBP (illustrated below), Compounds 501 to 509 below, or the like may be used.

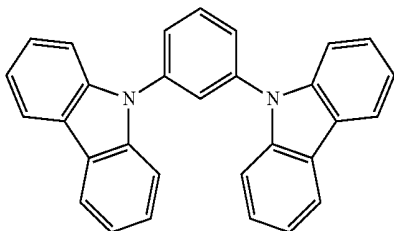
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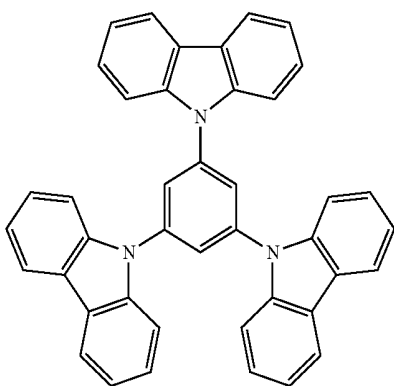
E3



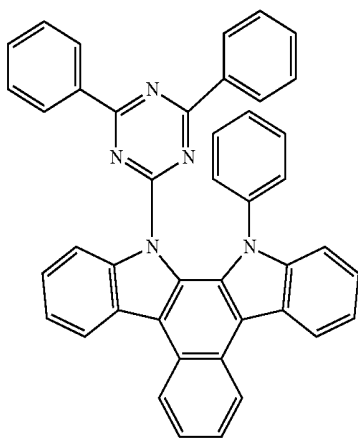
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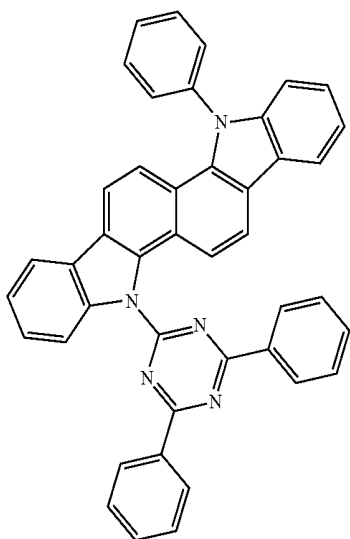
501



502

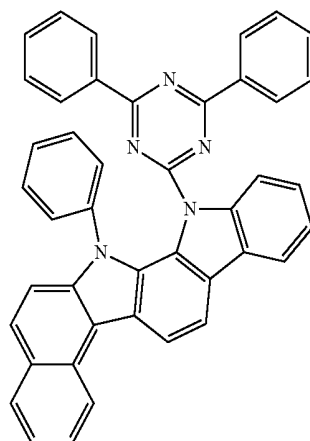


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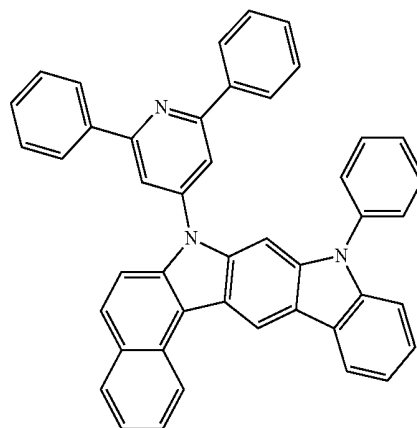


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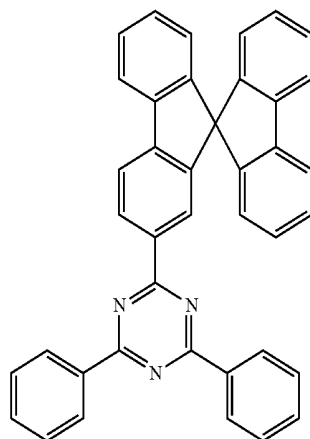
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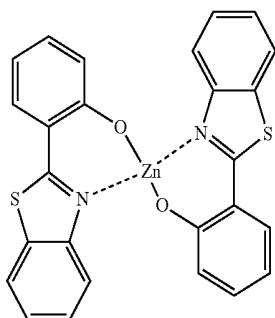


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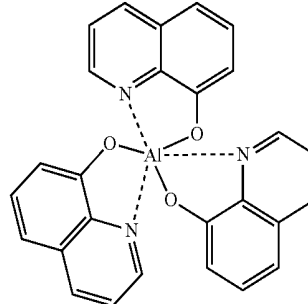
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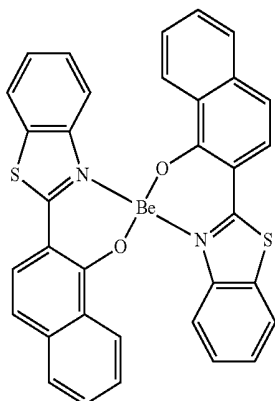


508

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Alq₃

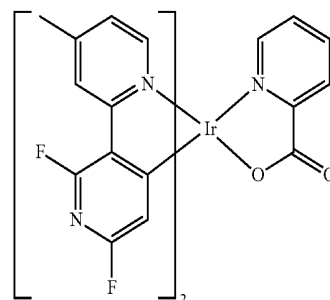
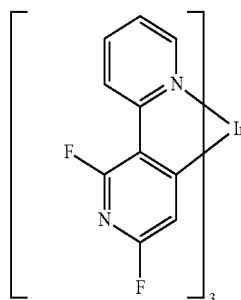
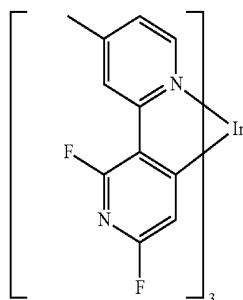
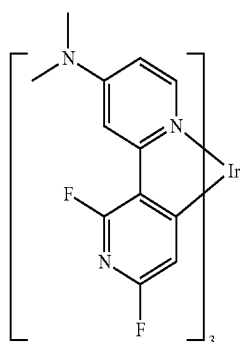
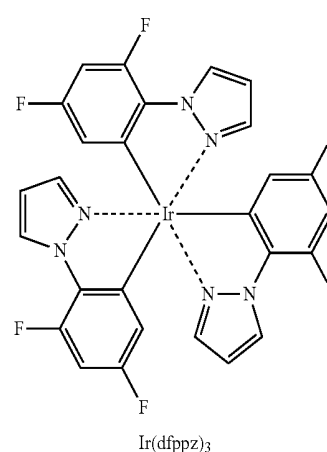
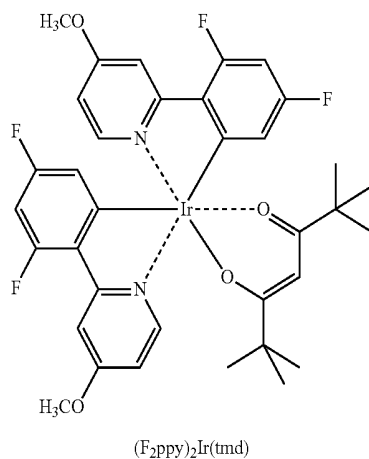
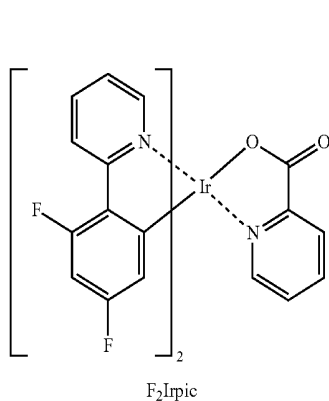
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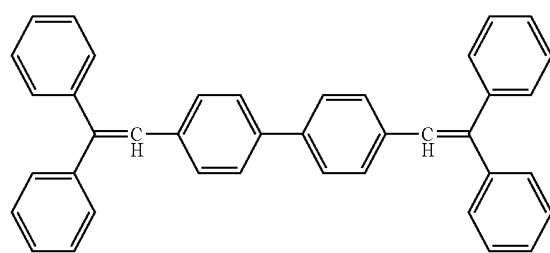
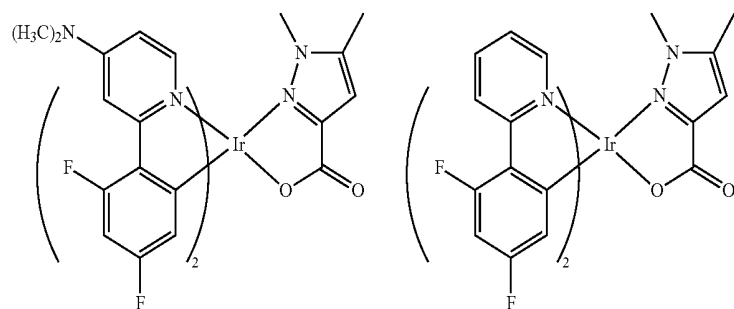
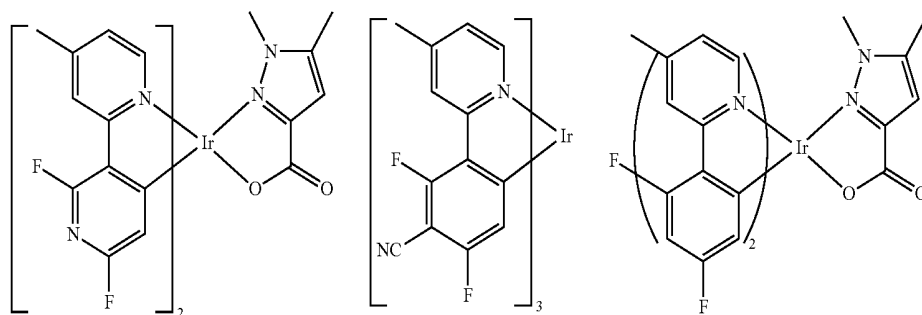
[0156] Alternatively, the EML may include as the host at least one compound selected from an anthracene-based compound represented by Formula 400 below and an anthracene-based compound represented by Formula 401.

[0157] The organic light-emitting diode may include a blue sub-pixel that emits blue light, a green sub-pixel that emits green light, and a red sub-pixel that emits red light. The blue sub-pixel may include a blue EML that emits blue light, and the blue EML may include the pyrene-based compound represented by Formula 1. According to another embodiment, due to a stack structure including a red EML, a green EML, and/or a blue EML, the EML may emit white light.

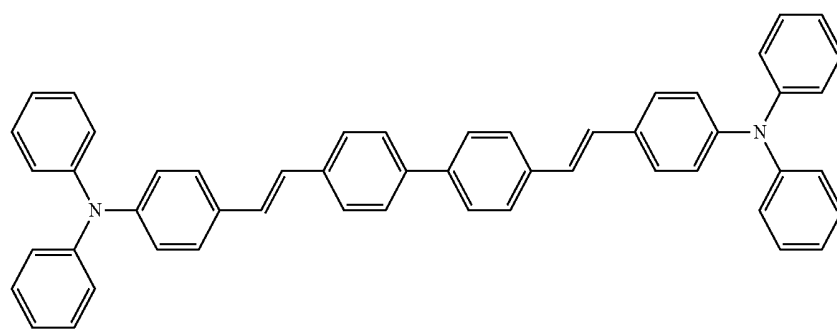
[0158] Also, the blue EML may further include compounds illustrated below as the blue dopant, but other materials may further be included in the blue EML.



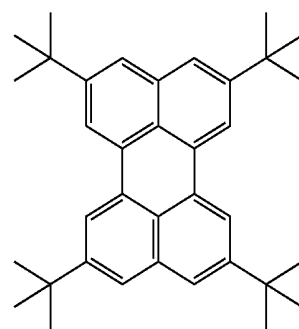
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DPVBi

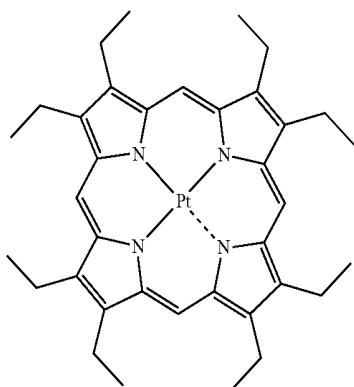


DPAVB

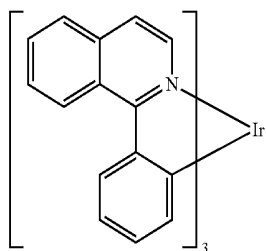
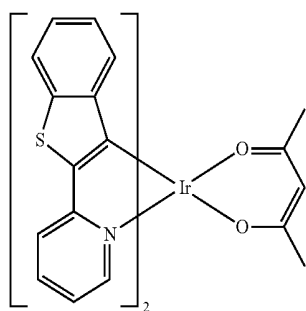
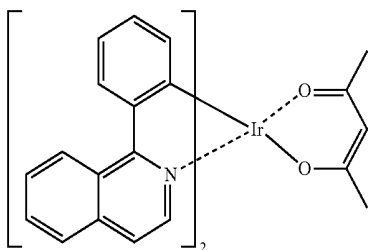


TBP

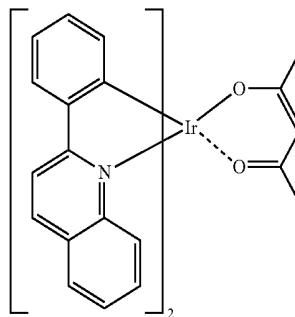
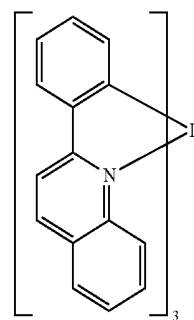
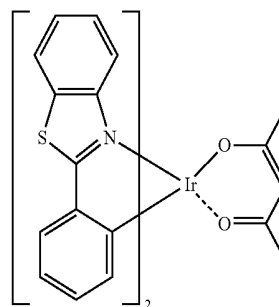
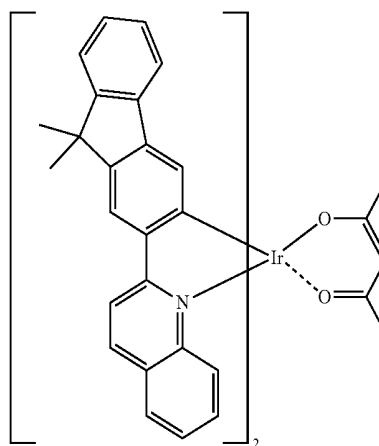
[0159] For example, the red EML of the red sub-pixel may include compounds illustrated below as a red dopant. In an implementation, DCM or DCJTb may be used as the red dopant.

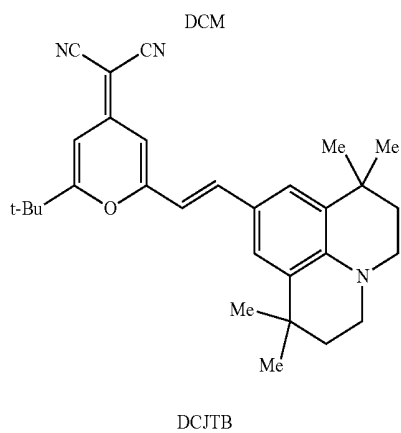
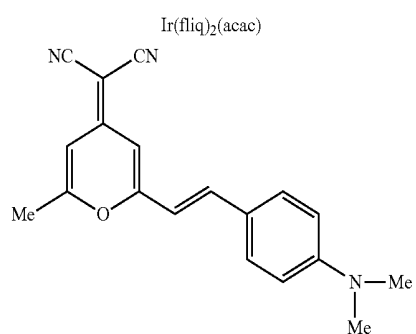
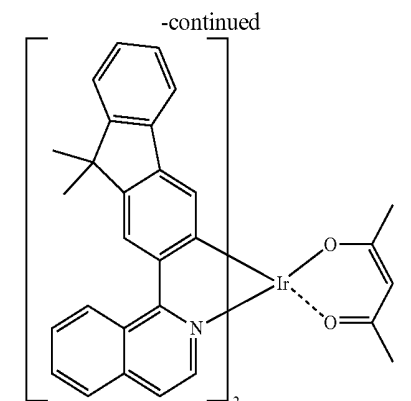


PtOEP

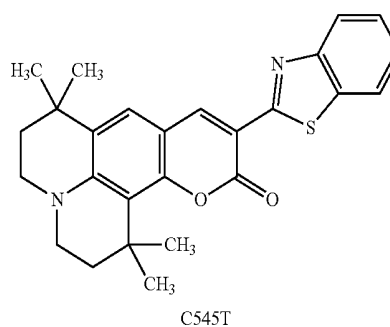
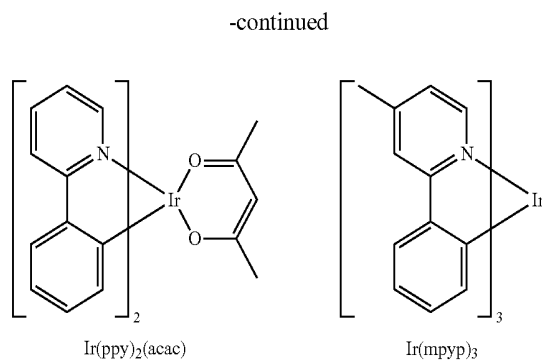
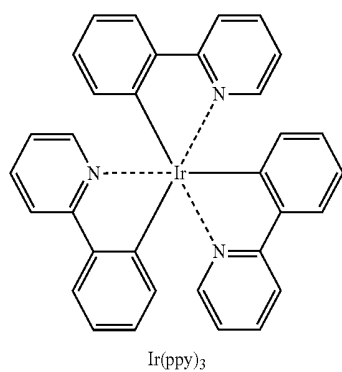
Ir(piq)₃Btp₂Ir(acac)

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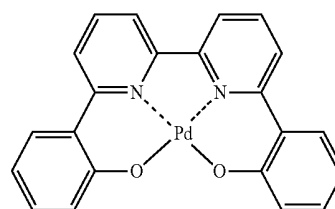
Ir(pq)₂(acac)Ir(2-phq)₃Ir(BT)₂(acac)Ir(flq)₂(acac)



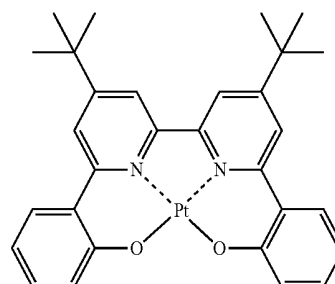
[0160] For example, the green EML of the green sub-pixel may include compounds illustrated below as a green dopant. In an implementation, a C545T below may be used as a green dopant.



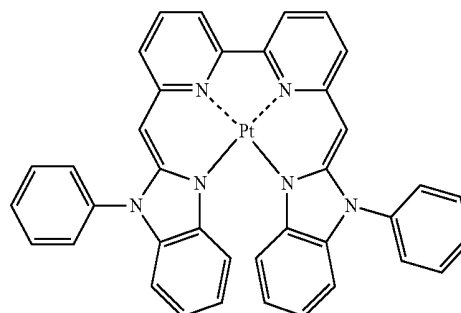
[0161] Another example of the dopant included in the EML may include a complex below.



D1

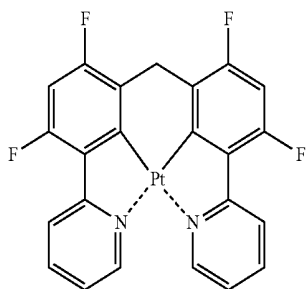
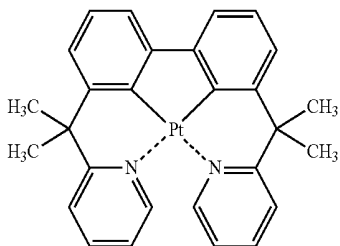
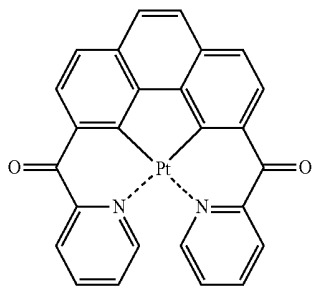
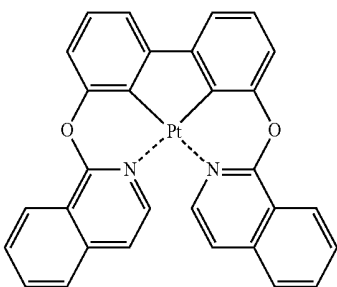
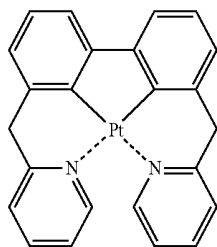
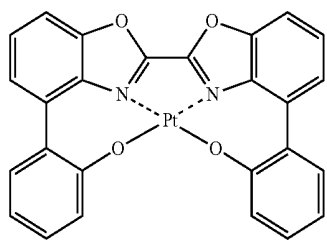


D2

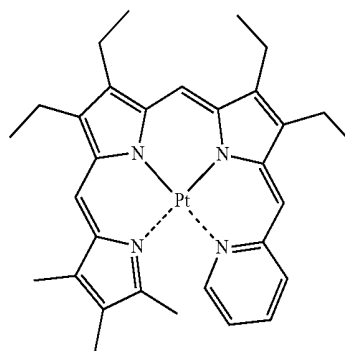
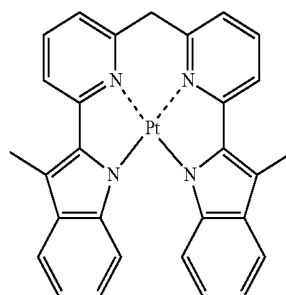
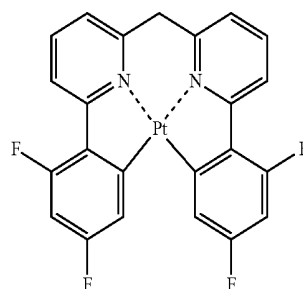
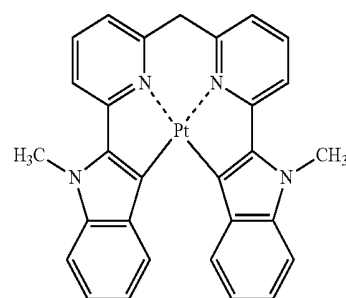
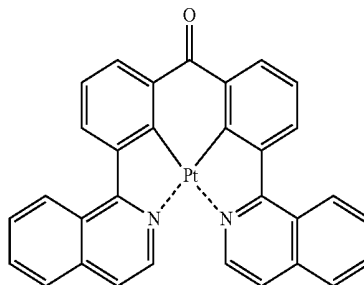


D3

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D4

D10

D5

D6

D7

D8

D9

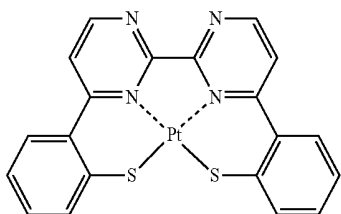
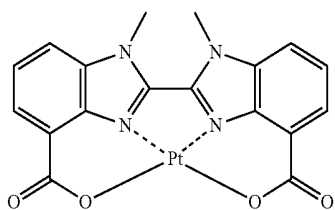
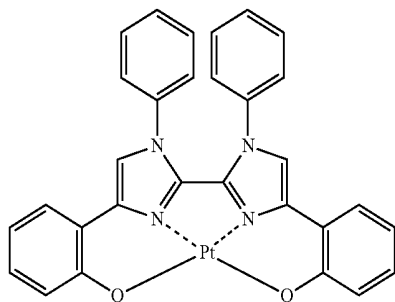
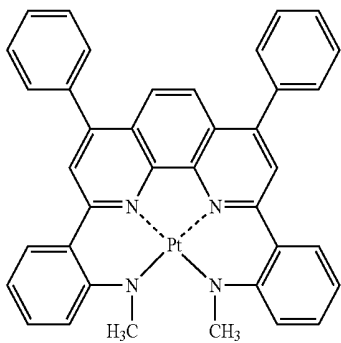
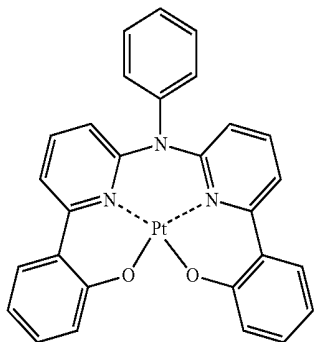
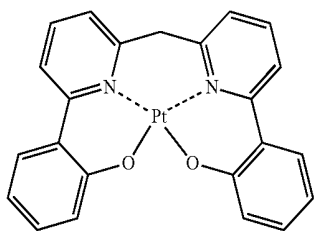
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D12

D13

D14

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D15

D16

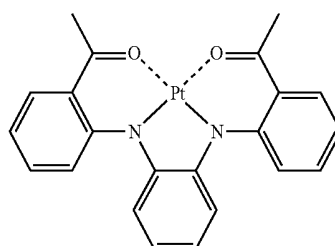
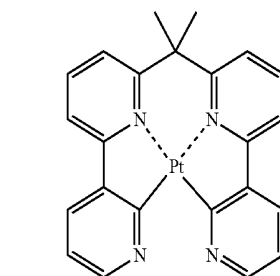
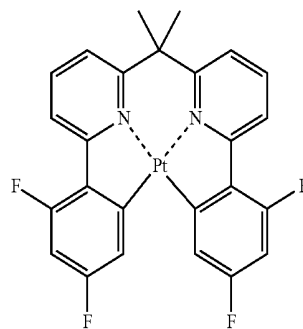
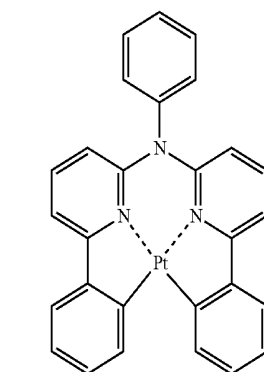
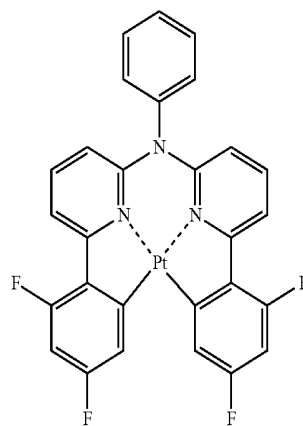
D17

D18

D19

D20

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D21

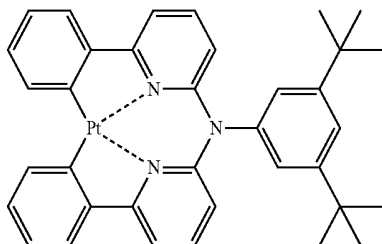
D22

D23

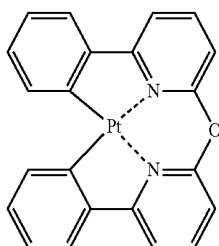
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D25

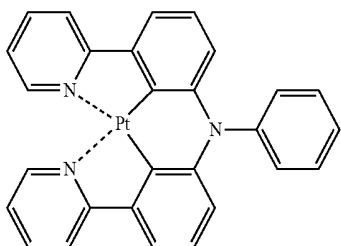
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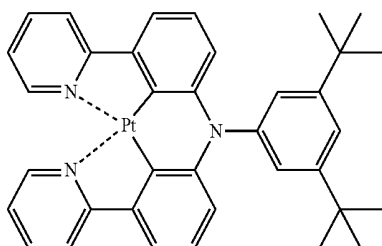
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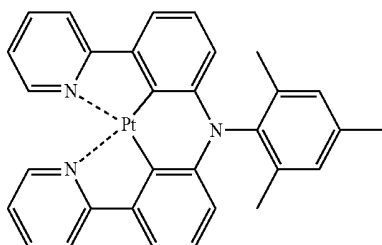
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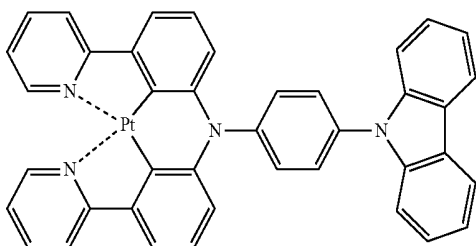
D28



D29

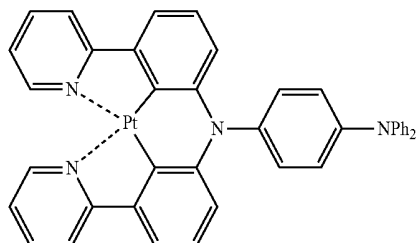


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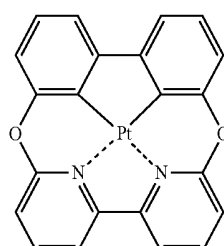


D31

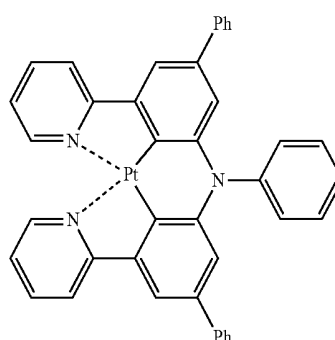
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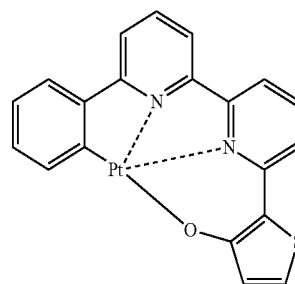
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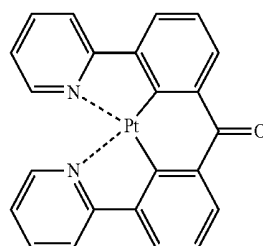
D33



D34

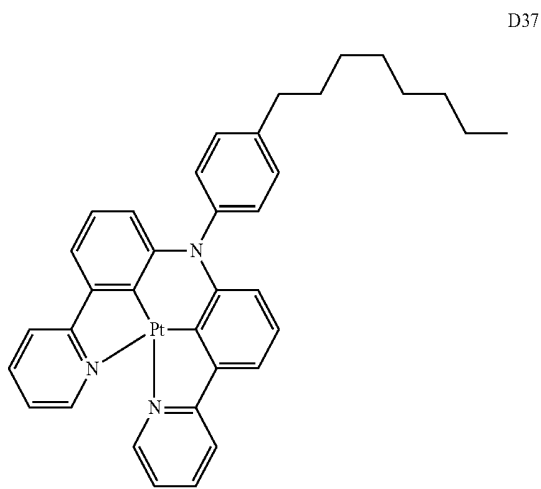


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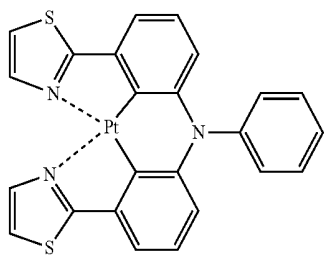


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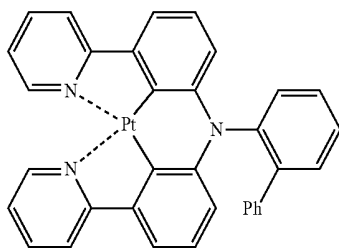
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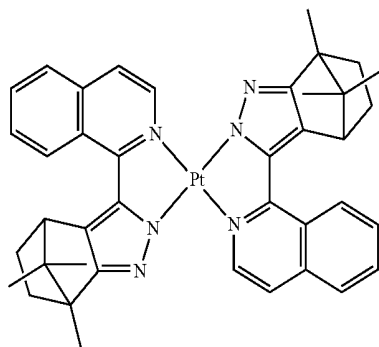
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D39

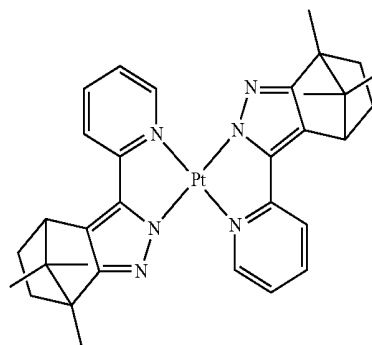


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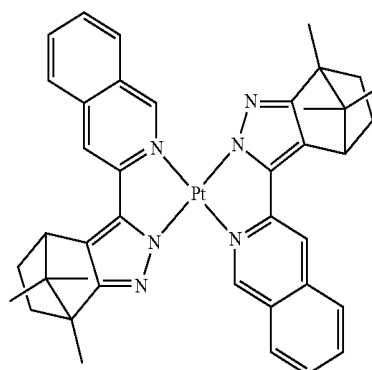


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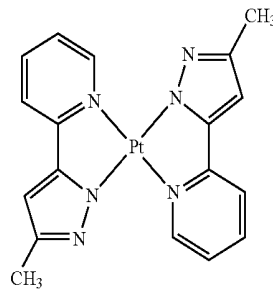
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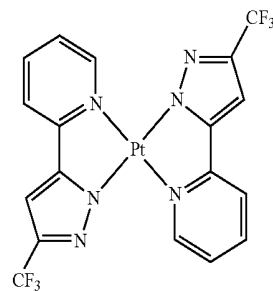
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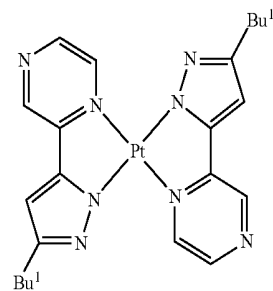
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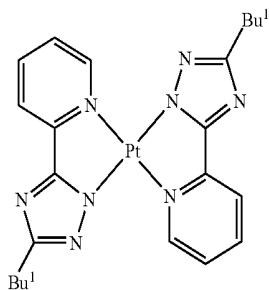
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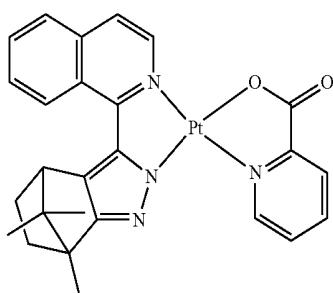
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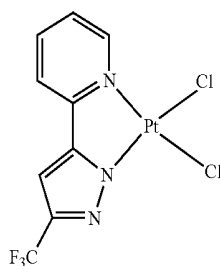
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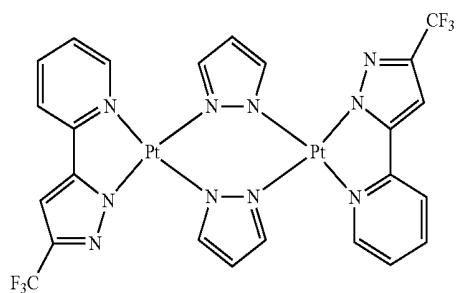
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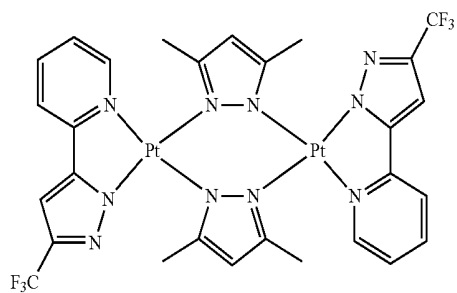
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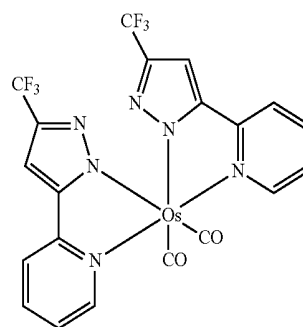
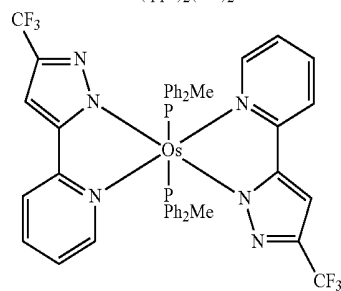
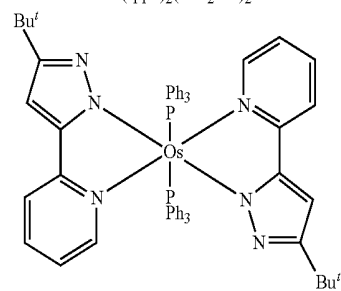
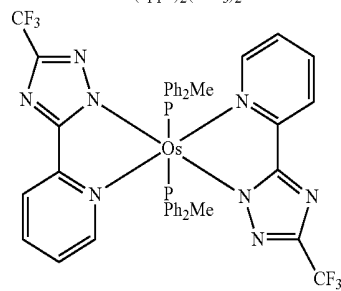
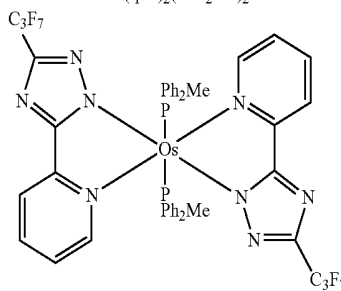
D48



D49



D50

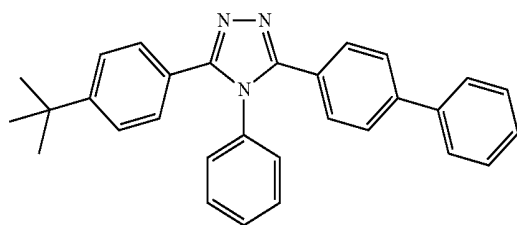
 $\text{Os}(\text{fppz})_2(\text{CO})_2$  $\text{Os}(\text{fppz})_2(\text{PPh}_2\text{Me})_2$  $\text{Os}(\text{bppz})_2(\text{PPh}_3)_2$  $\text{Os}(\text{fptz})_2(\text{PPh}_2\text{Me})_2$  $\text{Os}(\text{hptz})_2(\text{PPhMe}_2)_2$

[0162] Another example of the dopant included in the EML may include a Os-complex below.

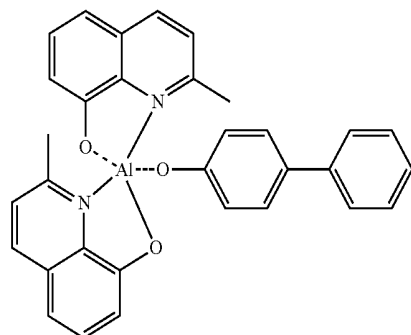
[0163] When the EML includes both a host and a dopant, an amount of the dopant may be from about 0.01 parts to about 15 parts by weight, based on 100 parts by weight of the host.

[0164] A thickness of the EML may be in a range of about 100 Å to about 1,000 Å, e.g., about 100 Å to about 600 Å. When the thickness of the EML is within these ranges, the EML may have improved luminescent ability without a substantial increase in driving voltage.

[0165] Next, an electron transport layer (ETL) may be formed on the EML using a suitable method, e.g., 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 conditions for deposition and coating may vary according to the material that is used to form the ETL. A material for an ETL may include suitable electron transporting materials that stably transport electrons injected from an electron injection electrode (cathode). Examples of the material for the ETL may include a quinoline derivative, such as tris(8-quinolinolate)aluminium (Alq3), TAZ, Balq, beryllium bis(benzoquinolin-10-olate) (Bebq2), ADN, Compound 201, and Compound 202.

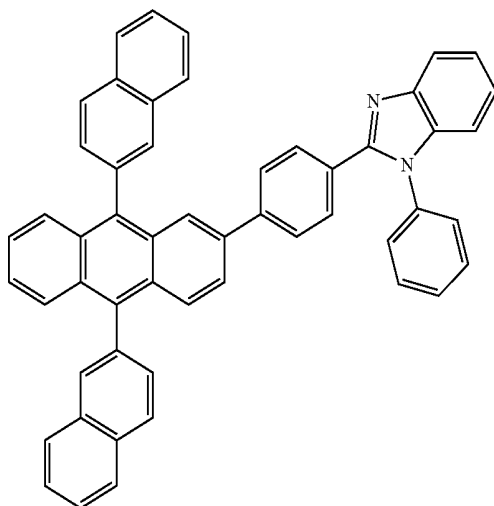


TAZ

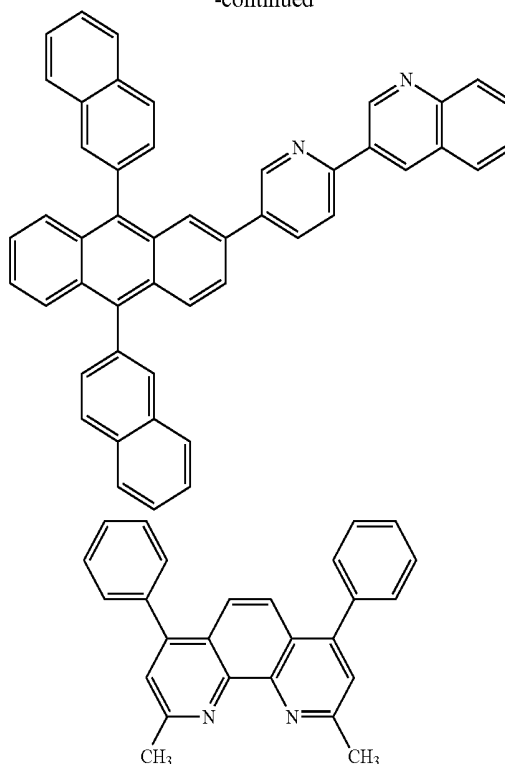


BAIq

<Compound 201>



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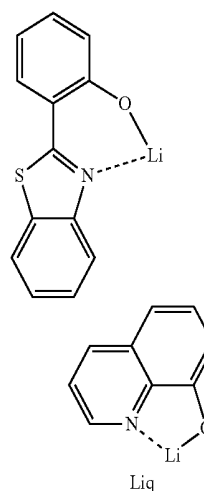


BCP

[0166] A thickness of the ETL may be in a range of about 100 Å to about 1,000 Å, e.g., about 150 Å to about 500 Å. When the thickness of the ETL is within these ranges, the ETL may have satisfactory electron transporting ability without a substantial increase in driving voltage.

[0167] The ETL may further include a metal-containing material, in addition to a suitable electron transporting inorganic material.

[0168] The metal-containing material may include a lithium (Li) complex. Examples of the Li complex may include lithium quinolate (Liq) and Compound 203 below.



Compound 203

Liq

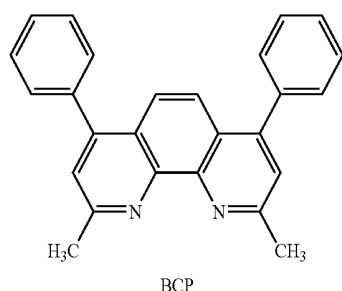
[0169] Then, an EIL, which facilitates injection of electrons from the cathode, may be formed on the ETL. A suitable electron-injecting material may be used to form the EIL.

[0170] Examples of materials for forming the EIL may include LiF, NaCl, a CsF, Li₂O, and BaO. The deposition conditions of the EIL may be similar to those used to form the HIL, although the deposition conditions may vary according to the material that is used to form the EIL.

[0171] The thickness of the EIL may be from about 1 Å to about 100 Å, e.g., may be from about 3 Å to about 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.

[0172] A second electrode 17 may be disposed on the organic layer 15. The second electrode 17 may be a cathode, which is an electron injecting electrode. A metal for forming the second electrode may be a metal, an alloy, an electrically conductive compound, which have a low-work function, or a mixture thereof. For example, lithium (Li), magnesium (Mg), aluminum (Al), aluminum (Al)-lithium (Li), calcium (Ca), magnesium (Mg)-indium (In), magnesium (Mg)-silver (Ag), or the like may be formed as a thin film to obtain a thin film-type reflective, semi-reflective, or a transmissive electrode. In an implementation, to manufacture a top-emission light-emitting device, a transmissive electrode formed of indium tin oxide (ITO) or indium zinc oxide (IZO) may be used. When the organic light-emitting diode is used in a large full-color display device, the second electrode (cathode) may be a reflective electrode.

[0173] In addition, when a phosphorescent dopant is used in the EML, a triplet exciton or a hole may diffuse to an ETL. To prevent the diffusion, a hole blocking layer (HBL) may be formed between the HTL and the EML or between the H-functional layer and the EML by vacuum deposition, spin coating, casting, LB deposition, or the like. When the HBL 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 HBL. A suitable hole-blocking material may be used. Examples of hole-blocking materials may include oxadiazole derivatives, triazole derivatives, and phenanthroline derivatives. For example, BCP illustrated below may be used as the hole-blocking material.



[0174] A thickness of the HBL may be in a range of about 20 Å to about 1,000 Å, e.g., about 30 Å to about 300 Å. When the thickness of the HBL is within these ranges, the HBL may have improved hole blocking ability without a substantial increase in driving voltage.

[0175] The organic light-emitting diode may be used in a full color display apparatus, a lamp, or the like. For example, the organic light-emitting diode may be a full color display apparatus.

[0176] Accordingly, provided is an organic light-emitting apparatus according to an embodiment including: a substrate including a first sub-pixel, a second sub-pixel, and a third sub-pixel; a first electrode formed in each of the first sub-pixel, second sub-pixel and third sub-pixel of the substrate; a second electrode that commonly faces the first electrodes in the first sub-pixel, the second sub-pixel, and the third sub-pixel; a first EML that is formed between the first electrode of the first sub-pixel and the second electrode and emits a first color light; a second EML that is formed between the first electrode of the second sub-pixel and the second electrode and emits a second color light; and a third EML that is formed between the first electrode of the third sub-pixel and the second electrode and emits a third color light, wherein the first EML includes one or more of the pyrene-based compounds described above. According to an embodiment, the first electrode may be a transmissive electrode or a semi-transmissive electrode, and the second electrode may be a reflective electrode. According to another embodiment, the first electrode may be a reflective electrode, and the second electrode may be a transmissive electrode or a semi-transmissive electrode.

[0177] A mixed light of the first light, the second light, and the third light of the organic light-emitting diode may be white light. Accordingly, the organic light-emitting diode may be a full-color display apparatus. The first light may be blue light. Also, the second light may be green light and the third light may be red light.

[0178] The first EML of the organic light-emitting apparatus may include the pyrene-based compound of Formula 1. Due to the inclusion of the pyrene-based compound, the emitted first color light (blue light) may have excellent color purity characteristics (e.g., a y color coordinate may be 1.0 or less) that satisfy NTSC or sRGB standards. Accordingly, the organic light-emitting apparatus may be used as a large high-quality TV.

[0179] According to an embodiment, the organic light-emitting apparatus may be a bottom emission type organic light-emitting apparatus in which the first electrode is a transmissive electrode or a semi-transmissive electrode, and the second electrode is a reflective electrode.

[0180] According to another embodiment, the organic light-emitting apparatus may be a top emission type organic light-emitting apparatus in which the first electrode is a reflective electrode, and the second electrode is a transmissive electrode or a semi-transmissive electrode.

[0181] The organic light-emitting apparatus may include the pyrene-based compound of Formula 1. Due to the inclusion of the pyrene-based compound, blue light having excellent color purity characteristics (e.g., a y color coordinate may be 1.0 or less) that satisfy a sRGB standard may be emitted. Accordingly, a complicated resonance structure that compensates for color purity of blue light is not needed, and thus, manufacturing costs may be reduced.

[0182] The full color display may be used in a television (TV), a personal computer monitor, a mobile communication terminal, an MP3 player, a navigation device for use in vehicles, or the like.

[0183] The unsubstituted C₁-C₆₀ alkyl group (or a C₁-C₆₀ alkyl group) used herein may be a C₁-C₆₀ linear or branched

alkyl group, such as methyl group, ethyl group, propyl group, isobutyl group, sec-butyl group, pentyl group, iso-amyl group, or hexyl group, and a substituent of a substituted C_1 - C_{60} alkyl group may be selected from a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine, a hydrazone, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkyne group, a C_1 - C_{60} alkoxy group; a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof; a C_3 - C_{10} cycloalkyl group, a C_3 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_3 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, and a C_2 - C_{60} heteroaryl group; a C_3 - C_{10} cycloalkyl group, a C_3 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_3 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, and a C_2 - C_{60} heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic 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 anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group; and $-N(Q_{11})(Q_{12})$; and $-Si(Q_{13})(Q_{14})(Q_{15})$ (in which Q_{11} and Q_{12} are each independently a C_6 - C_{60} aryl group, or a C_2 - C_{60} heteroaryl group, and Q_{13} to Q_{15} are each independently a C_1 - C_{60} alkyl group, a C_1 - C_{60} alkoxy group, a C_6 - C_{60} aryl group, or a C_2 - C_{60} heteroaryl group).

[0184] The unsubstituted C_1 - C_{60} alkoxy group (or C_1 - C_{60} alkoxy group) used herein may have a formula of $-OA$ (where A is the unsubstituted C_1 - C_{60} alkyl group described above), and detailed examples thereof are methoxy, ethoxy, and isopropoxy, and detailed examples of substituents of the unsubstituted C_1 - C_{60} alkoxy group may be understood by referring to the detailed examples of substituents of the substituted C_1 - C_{60} alkyl group.

[0185] The unsubstituted C_2 - C_{60} alkenyl group (or C_2 - C_{60} alkenyl group) used herein may refer to an unsubstituted C_2 - C_{60} alkyl group having one or more carbon double bonds at a center or end thereof. Examples of the unsubstituted C_2 - C_{60} alkenyl group are ethenyl group, propenyl group, and butenyl. Detailed examples of substituents of the substituted C_2 - C_{60} alkenyl groups may be understood by referring to the detailed examples of substituents of the substituted C_1 - C_{60} alkyl group.

[0186] The unsubstituted C_2 - C_{60} alkynyl group (or C_2 - C_{60} alkynyl group) used herein refers to an unsubstituted C_2 - C_{60} alkyl group having one or more carbon triple bonds at a center or end thereof. Examples of the unsubstituted C_2 - C_{60} alkynyl group are ethynyl group, propynyl group, and the like. Detailed examples of substituents of the substituted C_2 - C_{60}

alkynyl groups may be understood by referring to the detailed examples of substituents of the substituted C_1 - C_{60} alkyl group.

[0187] The unsubstituted C_3 - C_{10} cyclo alkyl group used herein may refer to a cyclic saturated hydrocarbon monovalent group, and examples thereof are cyclopropyl group, cyclobutyl group, cyclopentyl group, cyclohexyl group, cyclooctyl group, etc. Detailed examples of substituents of the substituted C_3 - C_{10} alkynyl groups may be understood by referring to the detailed examples of substituents of the substituted C_1 - C_{60} alkyl group.

[0188] The unsubstituted C_3 - C_{10} cycloalkenyl group may refer to a cyclic unsaturated hydrocarbon group that has one or more carbon double bonds and is not an aromatic cycle, and detailed examples thereof are cyclopropenyl group, cyclobutenyl group, cyclopentenyl group, cyclohexenyl group, cycloheptenyl group, 1,3-cyclohexadienyl group, 1,4-cyclohexadienyl group, 2,4-cycloheptadienyl group, and 1,5-cyclooctadienyl. Detailed examples of substituents of the substituted C_3 - C_{10} cycloalkenyl groups may be understood by referring to the detailed examples of substituents of the substituted C_1 - C_{60} alkyl group.

[0189] The unsubstituted C_6 - C_{60} aryl group may be a monovalent group having a carbocyclic aromatic system having 6 to 60 carbon atoms including at least one aromatic ring. The unsubstituted C_6 - C_{60} arylene group may be a divalent group having a carbocyclic aromatic system having 6 to 60 carbon atoms including at least one aromatic ring. When the aryl group and/or the arylene group have at least two rings, they may be fused to each other via a single bond. Detailed examples of substituents of the substituted C_6 - C_{60} aryl group and C_6 - C_{60} arylene group may be understood by referring to the detailed examples of substituents of the substituted C_1 - C_{60} alkyl group.

[0190] Examples of the substituted or unsubstituted C_6 - C_{60} aryl group may include a phenyl group, a C_1 - C_{10} alkylphenyl group (for example, ethylphenyl group), a C_1 - C_{10} alkylbiphenyl group (for example, ethylbiphenyl group), a halophenyl group (for example, an o-, m- or p-fluorophenyl group, a dichlorophenyl group), a dicyanophenyl group, a trifluoromethoxyphenyl group, o-, m-, and p-tolyl groups, o-, m- and p-cumenyl groups, 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, halonaphthyl group (for example, a fluoronaphthyl group), a C_1 - C_{10} alkyl naphthyl group (for example, a methyl naphthyl group), a C_1 - C_{10} alkoxynaphthyl group (for example, a methoxynaphthyl group), an anthracenyl group, azrenyl group, an acenaphthyl group, a phenalenyl group, a fluorenyl group, anthraquinolinyl group, a methylanthracenyl group, a phenanthrenyl group, a triphenylenyl group, a pyrenyl group, a chrysenyl group, an ethyl-chrysenyl group, a picenyl group, perylenyl group, a chloroperylene group, a pentaphenyl group, a pentasenyl group, a tetraphenylenyl group, a hexaphenyl group, a hexacenyl group, a rubicenyl group, a coronenyl group, a trinaphthyl group, a heptaphenyl group, a heptacenyl group, a piranthrenyl group, and an obarenyl group. Examples of the substituted or unsubstituted C_6 - C_{60} arylene group may be easily understood by referring to examples of the substituted or unsubstituted C_6 - C_{60} aryl group.

[0191] The unsubstituted C_2 - C_{60} heteroaryl group used herein may refer to a monovalent group having a system

composed of one or more aromatic rings having at least one hetero atom selected from nitrogen (N), oxygen (O), phosphorous (P), and sulfur (S) and carbon atoms as the remaining ring atoms. The unsubstituted C_2 - C_{60} heteroarylene group used herein may refer to a divalent group having a system composed of one or more aromatic rings having at least one hetero atom selected from nitrogen (N), oxygen (O), phosphorous (P), and sulfur (S) and carbon atoms as the remaining ring atoms. In this regard, when the heteroaryl group and the heteroarylene group each include two or more rings, the rings may be fused to each other. Detailed examples of substituents of the substituted C_2 - C_{60} heteroaryl group and C_2 - C_{60} heteroarylene group may be understood by referring to the detailed examples of substituents of the substituted C_1 - C_{60} alkyl group.

[0192] Examples of the unsubstituted C_2 - C_{60} heteroaryl group may include a pyrazolyl group, an imidazolyl group, a 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, benzoan imidazolyl group, an imidazo pyridinyl group, and an imidazo pyrimidinyl group. Examples of the substituted or unsubstituted C_2 - C_{60} heteroarylene group may be easily understood by referring to examples of the substituted or unsubstituted C_2 - C_{60} heteroaryl group.

[0193] The substituted or unsubstituted C_6 - C_{60} aryloxy group may be represented by $-OA_2$ (wherein A_2 indicates the substituted or unsubstituted C_6 - C_{60} aryl group), and the substituted or unsubstituted C_6 - C_{60} arylthio group may be represented by $-SA_3$ (wherein A_3 indicates a substituted or unsubstituted C_6 - C_{60} aryl group).

[0194] The following Examples and Comparative Examples are provided in order to highlight characteristics of one or more embodiments, but it will be understood that the Examples and Comparative Examples are not to be construed as limiting the scope of the embodiments, nor are the Comparative Examples to be construed as being outside the scope of the embodiments. Further, it will be understood that the embodiments are not limited to the particular details described in the Examples and Comparative Examples.

EXAMPLE

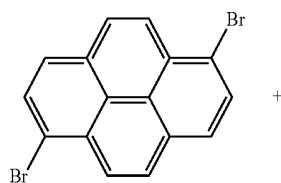
[0195] Intermediates 1 to 14 were synthesized by using methods disclosed in i) Journal of Medicinal Chemistry, 51(23), 7640-7644, 2008, ii) Synthesis (5), 815-823, 2009, and iii) Bioorganic & Medicinal Chemistry 19(2), 939-950, 2011, which are incorporated by references herein in their entirety.

Synthesis Example 1

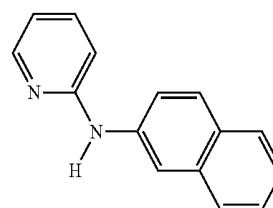
Synthesis of Compound 4

Synthesis of Intermediate PY1

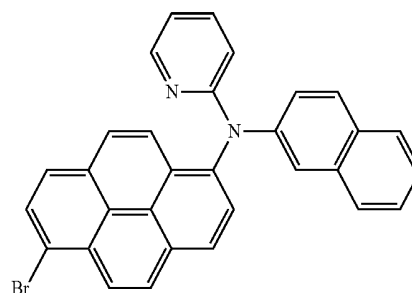
[0196]



-continued



Intermediate 1

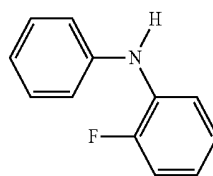


Intermediate PY1

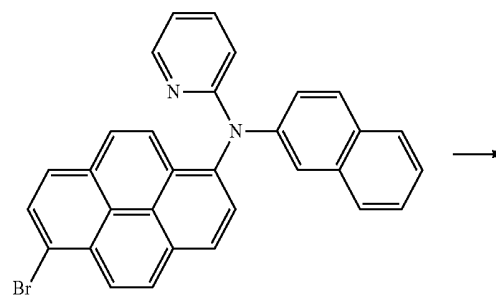
[0197] Under nitrogen atmosphere, 3.6 g (10 mmol) of 1,6-dibromopyrene, 2.2 g (10 mmol) of Intermediate 1, 287 mg (0.5 mmol) of bis(dibenzylideneacetone)palladium ($Pd(dba)_2$), 303 mg (1.5 mmol) of $P(tert-Bu)_3$, and 2.9 g (30 mmol) of sodium t-butoxide were added to 500 ml of toluene, and then, the mixture was refluxed for 12 hours. When the reaction was completed, a solvent was removed by evaporation, and then, 500 ml of methylene chloride and 500 ml of water were respectively added thereto and the mixture was washed to collect an organic layer which was then dried with an anhydrous magnesium sulfate. Subsequently, the result was subjected to recrystallization and silica gel chromatography to obtain Intermediate PY1 (1 g, yield of 20%).

Synthesis of Compound 1

[0198]

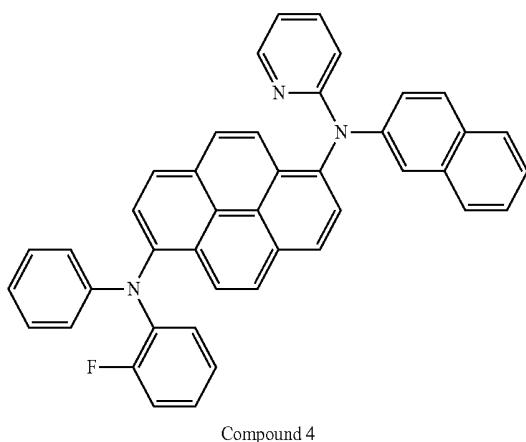


Intermediate 2



Intermediate PY1

-continued



[0199] Compound 4 (303 mg, yield of 49%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 2 and Intermediate PY1 were respectively used instead of 1,6-dibromopyrene and Intermediate 1.

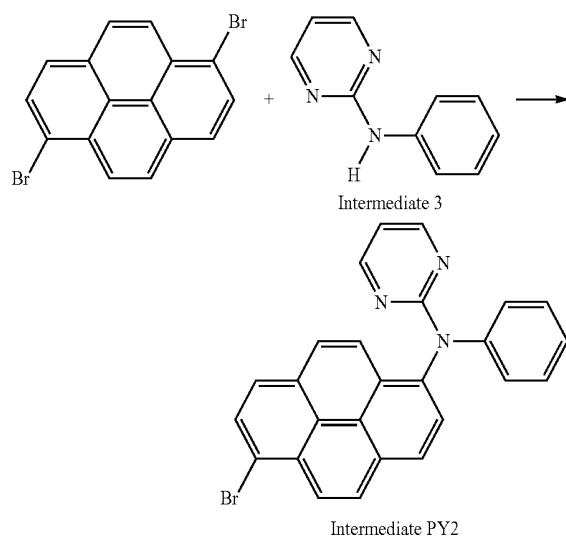
[0200] MS (MALDI-TOF) m/z : 605 $[M]^+$.

Synthesis Example 2

Synthesis of Compound 7

Synthesis of Intermediate PY2

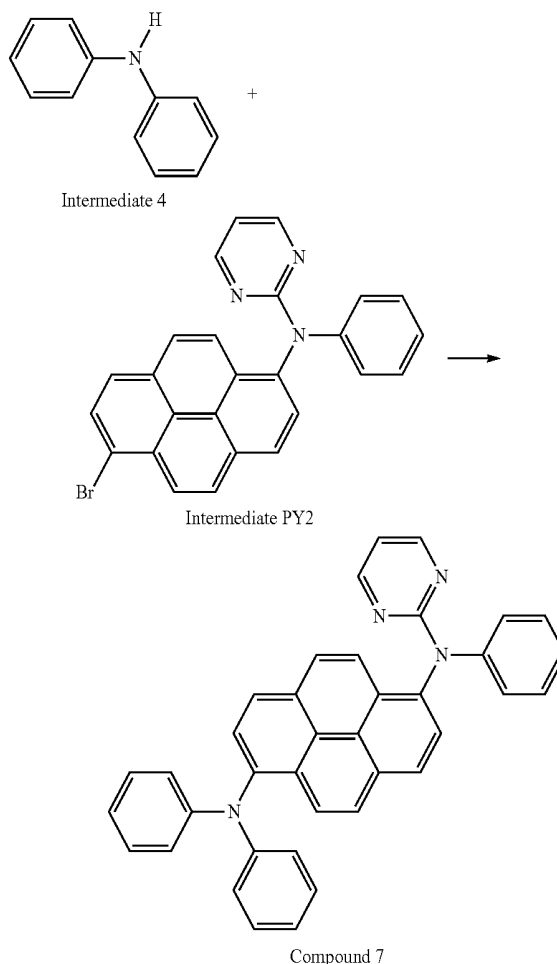
[0201]



[0202] Intermediate PY2 (502 mg, yield of 12%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 3 was used instead of Intermediate 1.

Synthesis of Compound 1

[0203]



[0204] Compound 7 (298 mg, yield of 48%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 4 and Intermediate PY2 were respectively used instead of 1,6-dibromopyrene and Intermediate 1.

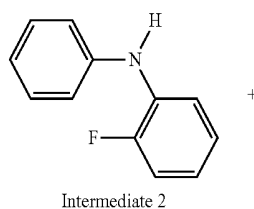
[0205] MS (MALDI-TOF) m/z : 538 $[M]^+$.

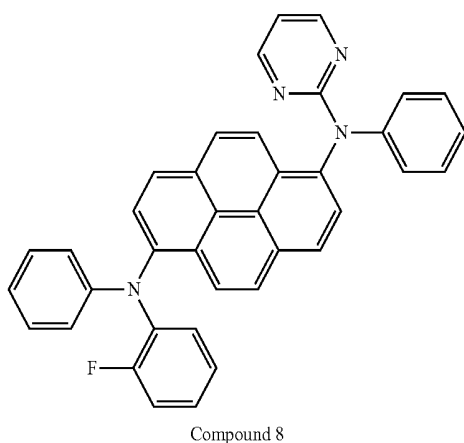
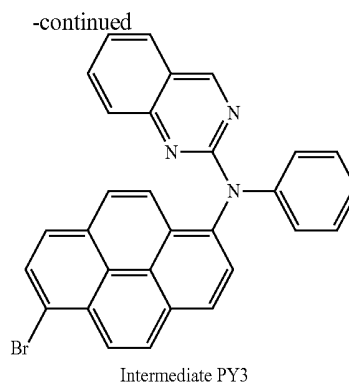
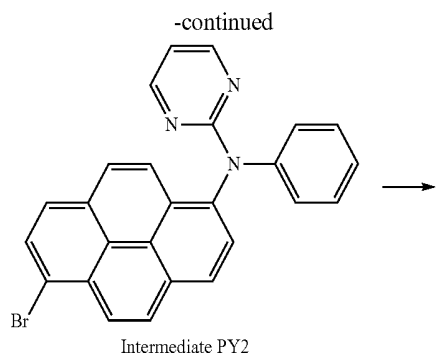
Synthesis Example 3

Synthesis of Compound 8

Synthesis of Compound 8

[0206]





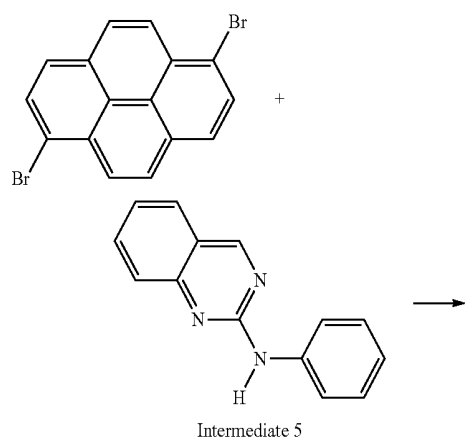
[0207] Compound 8 (338 mg, yield of 43%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 2 and Intermediate PY2 were respectively used instead of 1,6-dibromopyrene and Intermediate 1. MS (MALDI-TOE) m/z : 556 $[M]^+$.

Synthesis Example 4

Synthesis of Compound 13

Synthesis of Intermediate PY3

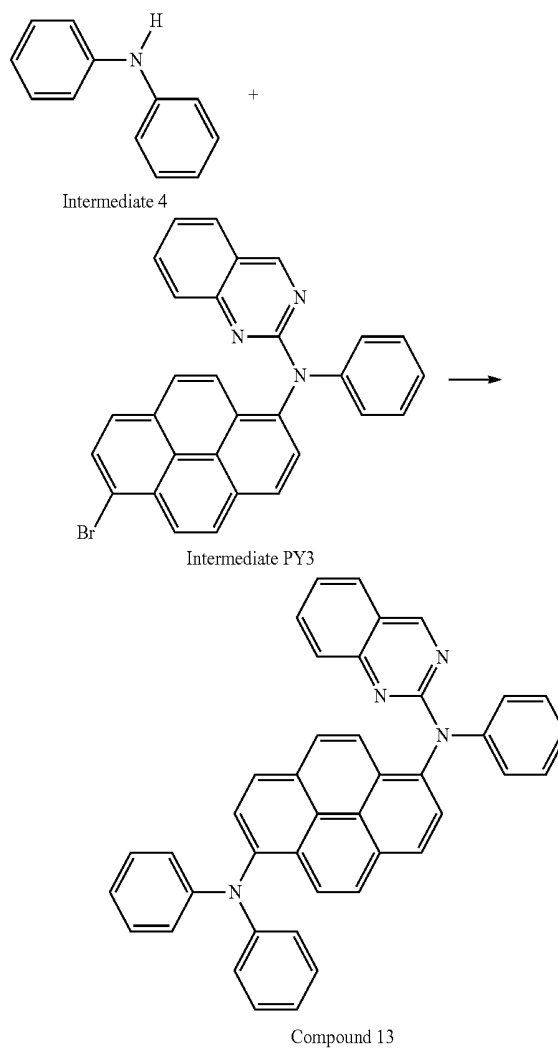
[0208]



[0209] Intermediate PY3 (278 mg, yield of 10%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 5 was used instead of Intermediate 1.

Synthesis of Compound 13

[0210]



[0211] Compound 13 (502 mg, yield of 44%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 4 and Intermediate PY3 were respectively used instead of 1,6-dibromopyrene and Intermediate 1.

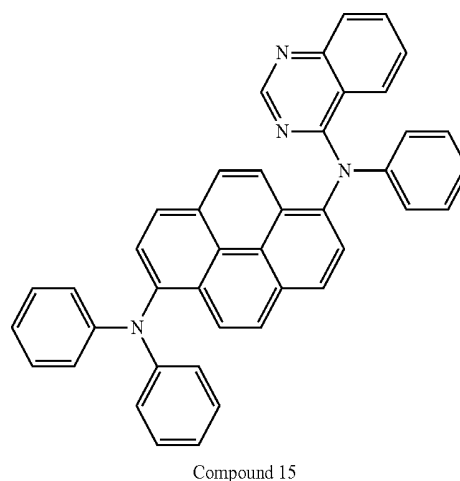
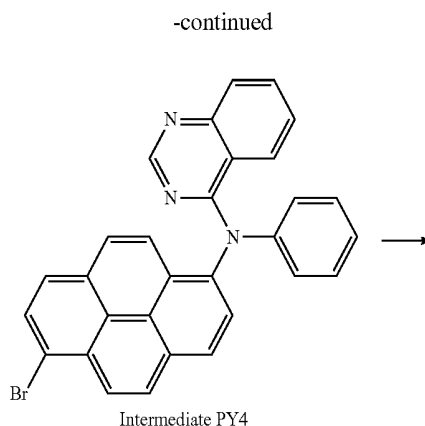
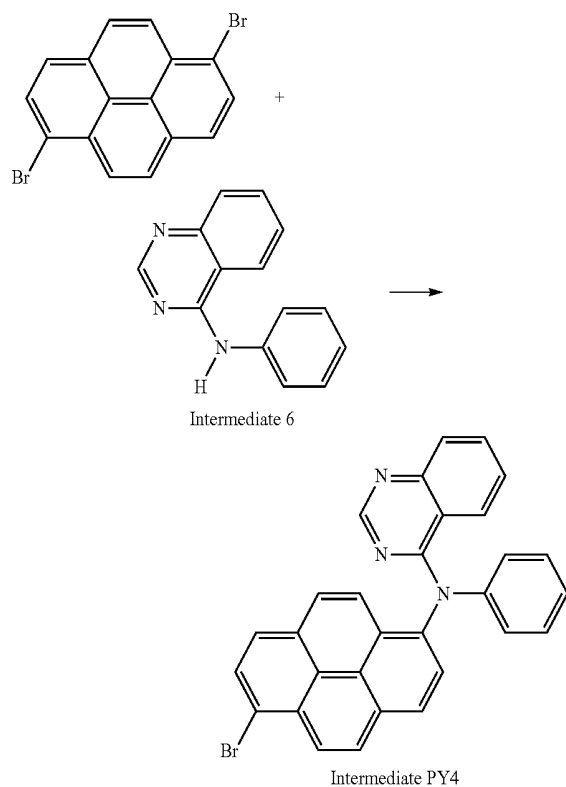
[0212] MS (MALDI-TOF) m/z : 588 $[M]^+$.

Synthesis Example 5

Synthesis of Compound 15

Synthesis of Intermediate PY4

[0213]



[0216] Compound 15 (679 mg, yield of 35%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 4 and Intermediate PY4 were respectively used instead of 1,6-dibromopyrene and Intermediate 1.

[0217] MS (MALDI-TOF) m/z : 588 $[M]^+$.

Synthesis Example 6

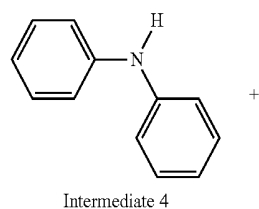
Synthesis of Compound 32

Synthesis of Intermediate PY5

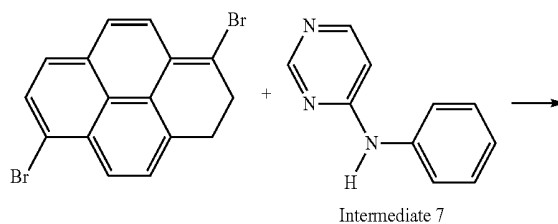
[0214] Intermediate PY4 (444 mg, yield of 13%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 6 was used instead of Intermediate 1.

Synthesis of Compound 15

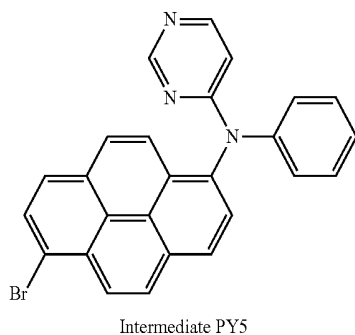
[0215]



[0218]



-continued

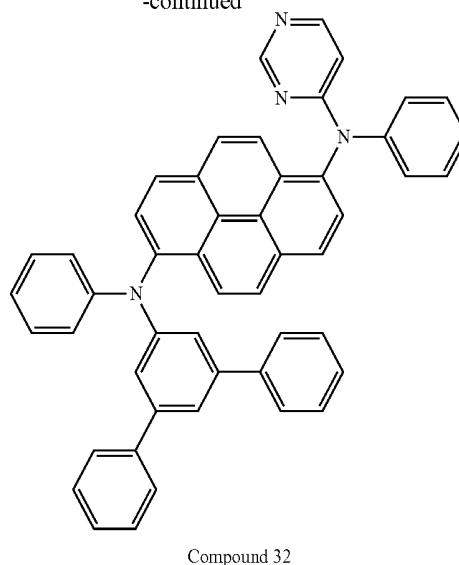


[0219] Intermediate PY5 (275 mg, yield of 12%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 7 was used instead of Intermediate 1.

Synthesis of Compound 32

[0220]

-continued



[0221] Compound 32 (459 mg, yield of 39%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 8 and Intermediate PY5 were respectively used instead of 1,6-dibromopyrene and Intermediate 1.

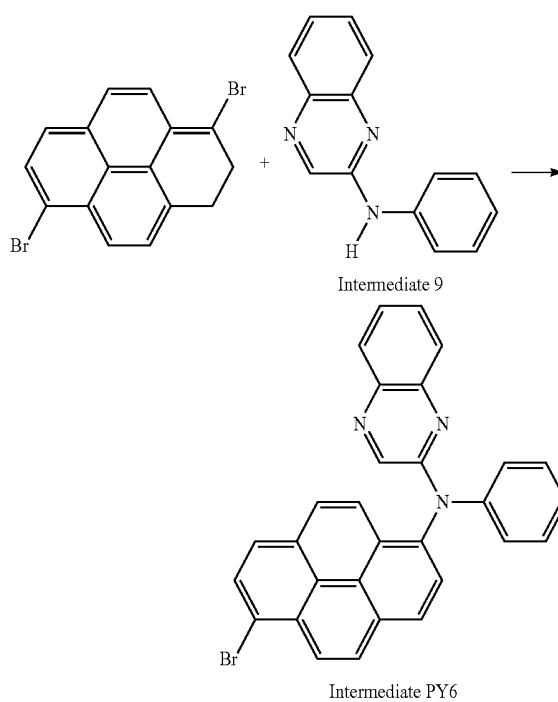
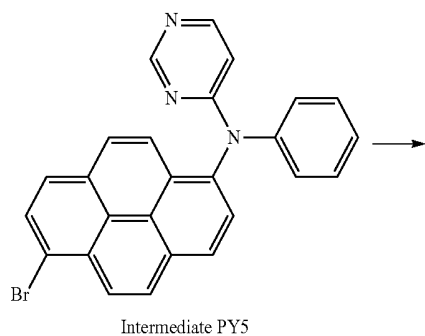
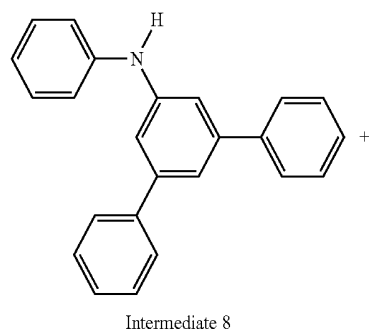
[0222] MS (MALDI-TOF) m/z: 690 [M]⁺.

Synthesis Example 7

Synthesis of Compound 35

Synthesis of Intermediate PY6

[0223]



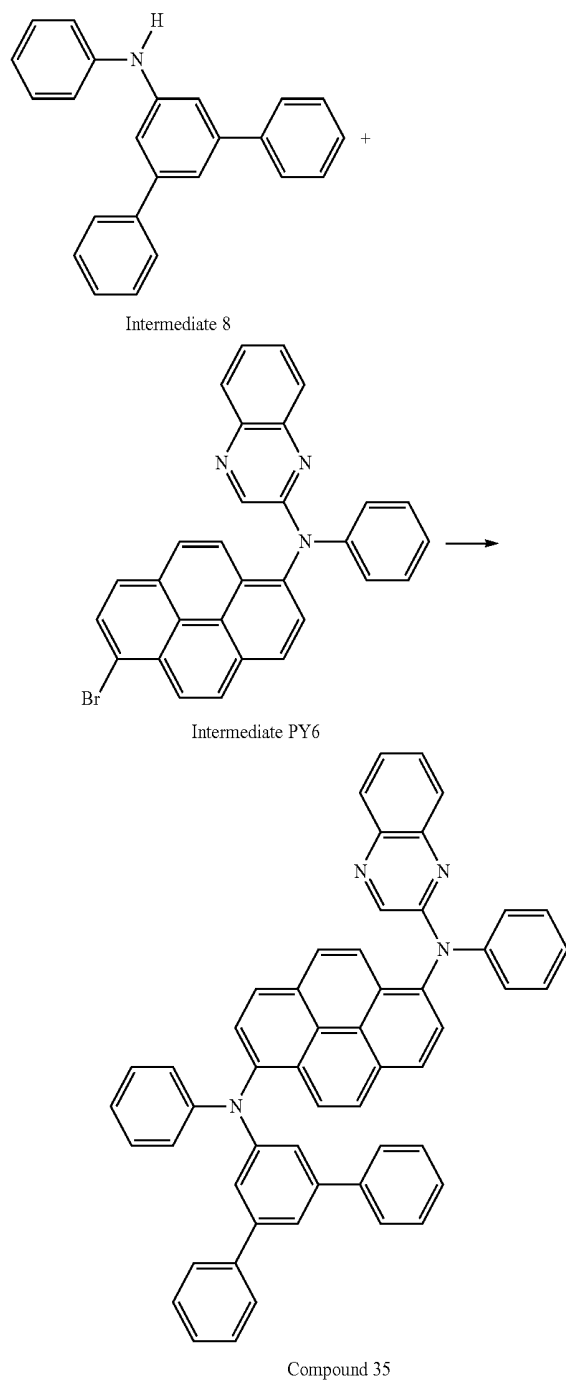
[0224] Intermediate PY6 (351 mg, yield of 13%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 9 was used instead of Intermediate 1.

respectively used instead of 1,6-dibromopyrene and Intermediate 1.

[0227] MS (MALDI-TOF) m/z : 740 $[M]^+$.

Synthesis of Compound 35

[0225]



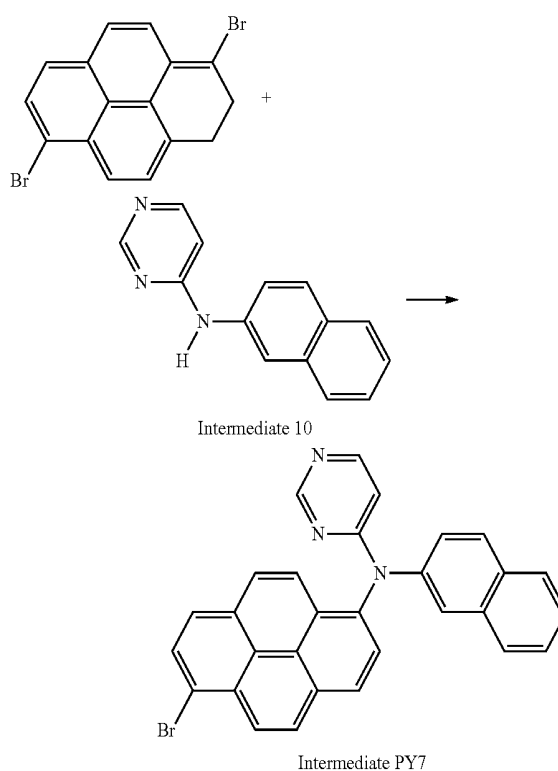
[0226] Compound 35 (672 mg, yield of 46%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 8 and Intermediate PY6 were

Synthesis Example 8

Synthesis of Compound 44

Synthesis of Intermediate PY7

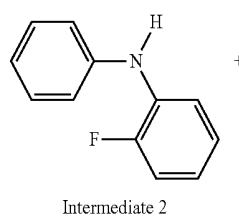
[0228]



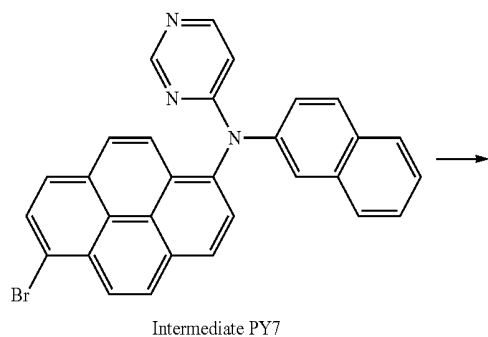
[0229] Intermediate PY7 (358 mg, yield of 11%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 10 was used instead of Intermediate 1.

Synthesis of Compound 44

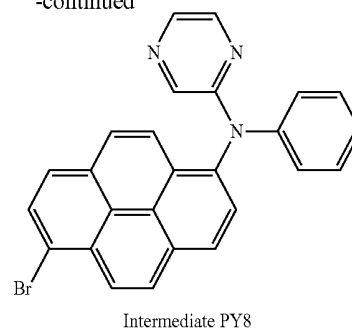
[0230]



-continued



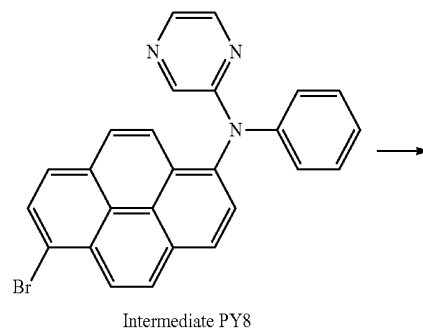
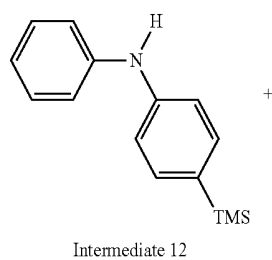
-continued



[0234] Intermediate PY8 (498 mg, yield of 12%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 11 was used instead of Intermediate 1.

Synthesis of Compound 69

[0235]



[0231] Compound 44 (521 mg, yield of 43%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 2 and Intermediate PY7 were respectively used instead of 1,6-dibromopyrene and Intermediate 1.

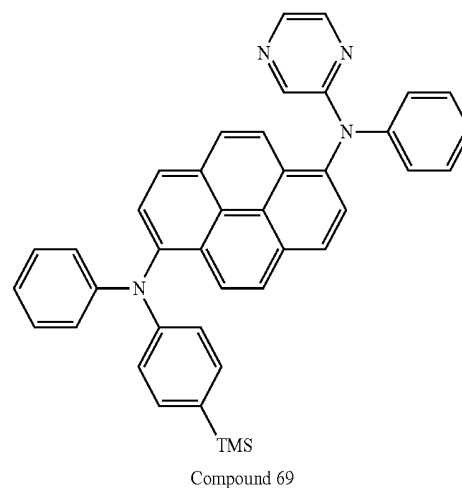
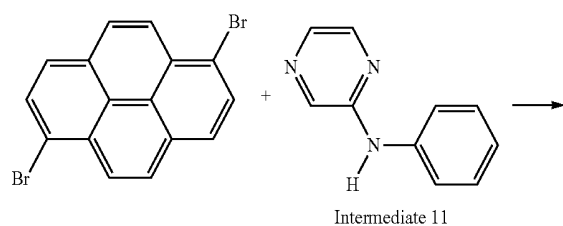
[0232] MS (MALDI-TOF) m/z : 606 $[M]^+$.

Synthesis Example 9

Synthesis of Compound 69

Synthesis of Intermediate PY8

[0233]



[0236] Compound 69 (475 mg, yield of 41%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 12 and Intermediate PY8 were respectively used instead of 1,6-dibromopyrene and Intermediate 1.

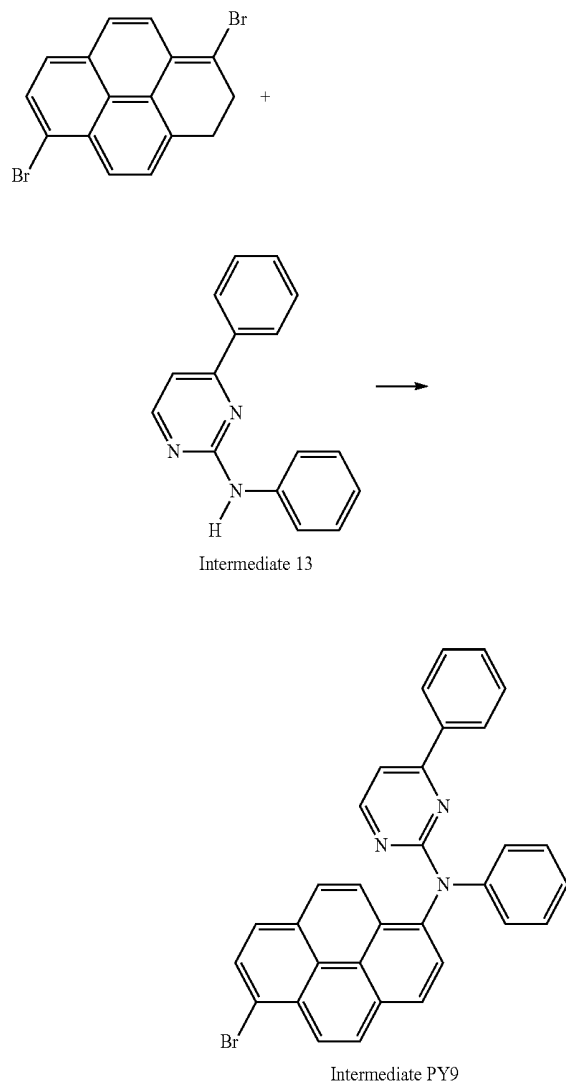
[0237] MS (MALDI-TOF) m/z : 610 $[M]^+$.

Synthesis Example 10

Synthesis of Compound 87

Synthesis of Intermediate PY9

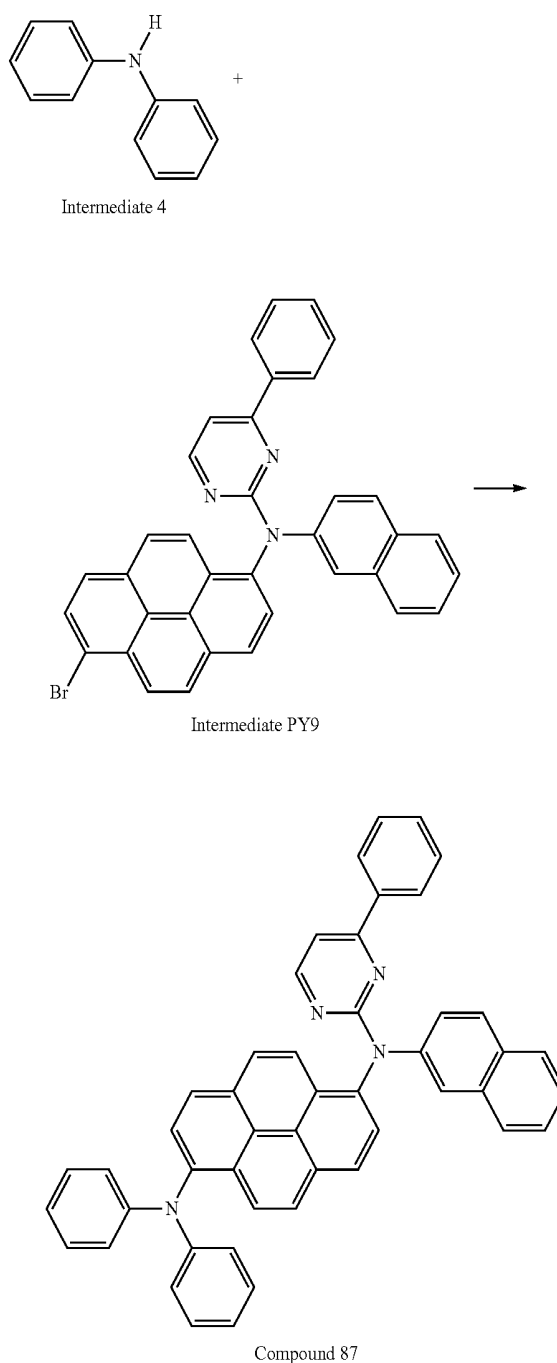
[0238]



[0239] Intermediate PY9 (277 mg, yield of 9%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 13 was used instead of Intermediate 1.

Synthesis of Compound 87

[0240]



[0241] Compound 87 (535 mg, yield of 47%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 4 and Intermediate PY9 were respectively used instead of 1,6-dibromopyrene and Intermediate 1.

[0242] MS (MALDI-TOF) m/z : 664 $[M]^+$.

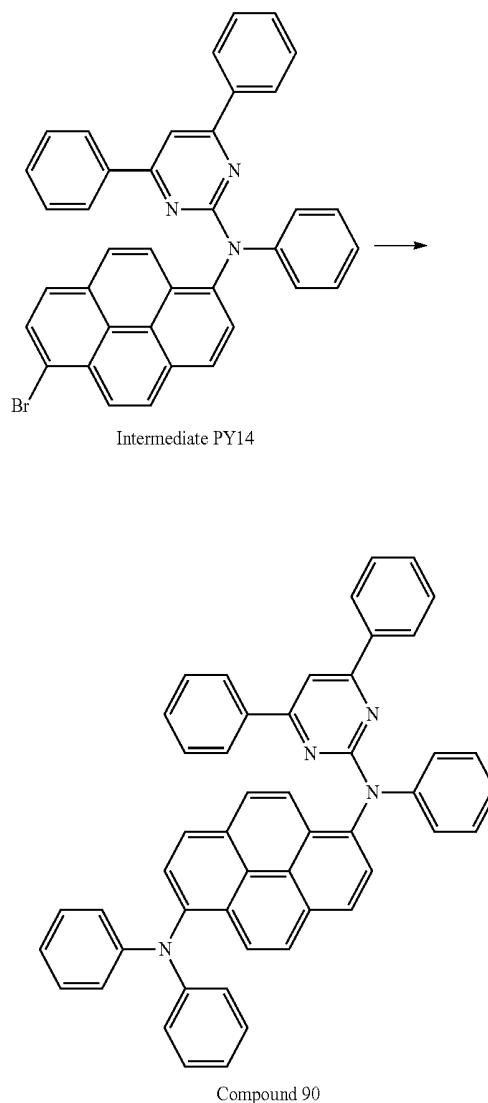
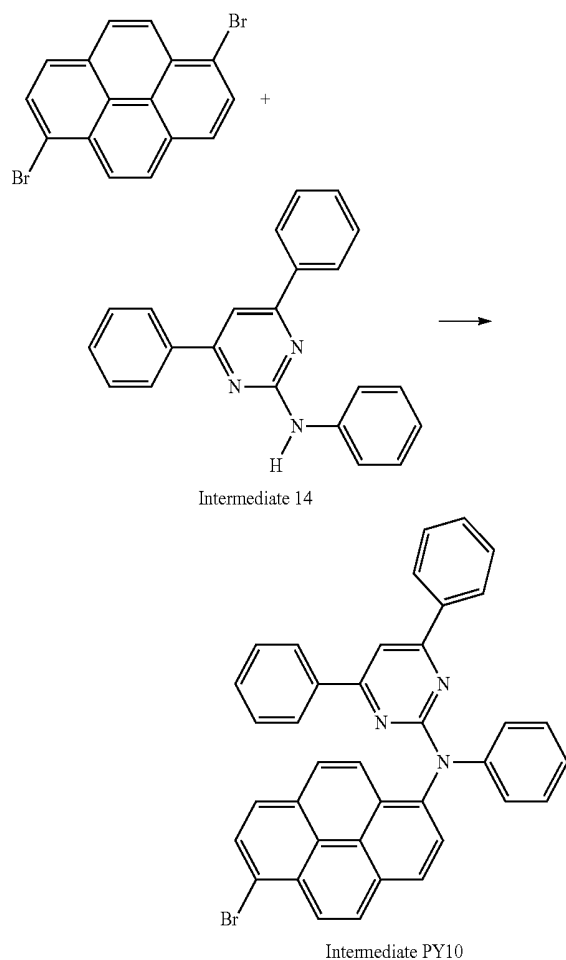
Synthesis Example 11

-continued

Synthesis of Compound 90

Synthesis of Intermediate PY10

[0243]



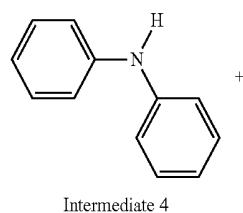
[0244] Intermediate PY10 (378 mg, yield of 13%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 14 was used instead of Intermediate 1.

[0246] Compound 90 (423 mg, yield of 41%) was synthesized in the same manner as used in synthesizing Intermediate PY1, except that Intermediate 4 and Intermediate PY14 were respectively used instead of 1,6-dibromopyrene and Intermediate 1.

[0247] MS (MALDI-TOF) m/z : 690 $[M]^+$.

Synthesis of Compound 90

[0245]



Evaluation Example 1

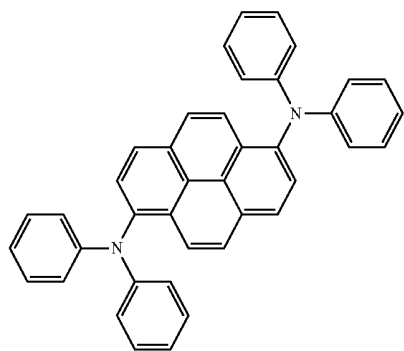
PL Spectrum Evaluation

[0248] Compound 7 was dissolved at a concentration of $1 \times 10^{-5} M$ in $CHCl_3$, and a photoluminescence (PL) spectrum thereof was measured by using ISC PC1 spectrofluorometer equipped with a Xenon lamp. The same experiment was performed with Compound 8 and Compounds A and B below to obtain PL spectrum maximum wavelengths, and results thereof are shown in Table 1, below.

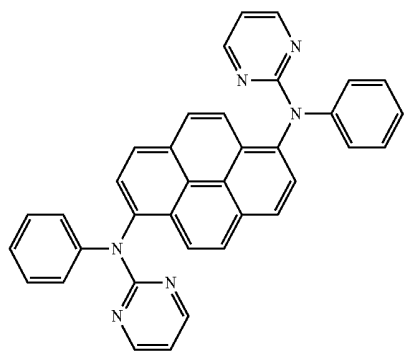
TABLE 1

	Compound A	Compound 7	Compound 8	Compound B
PL maximum wavelength (nm)	456	452	437	419

<Compound A>



<Compound B>



[0249] Referring to Table 1, it may be seen that Compounds 7 and 8 exhibited better blue color purity characteristics than Compound A, and Compound B exhibited PL characteristics with a relatively short wavelength region.

Example 1

[0250] A Corning 15 Ωcm^2 (1,200 Å) ITO glass substrate was cut to a size of 50 mm×50 mm×0.5 mm, and then, the ITO glass substrate was sonicated with isopropyl alcohol and pure water each for 5 minutes, and then cleaned by exposure to UV light for 30 minutes and to ozone. Then, the ITO glass substrate was mounted on a vacuum deposition apparatus.

[0251] m-MTDATA was deposited on the ITO glass substrate functioning as an anode at a deposition rate of 1 Å/sec to form a hole injection layer having a thickness of 600 Å, and then alpha-NPD was deposited on the hole injection layer at a deposition rate of 1 Å/sec to form a hole transport layer having a thickness of 300 Å.

[0252] Then, Compound 4(dopant) and 9,10-di-naphthalene-2-yl-anthracene (AND, host) were co-deposited on the hole transport layer at deposition rates of 0.05 Å/sec and 1 Å/sec, respectively, to form an emission layer having a thickness of 200 Å.

[0253] Thereafter, Alq₃ was deposited on the emission layer to form an electron transport layer having a thickness of 300 Å, and LiF was deposited on the electron transport layer to form an electron injection layer having a thickness of 10 Å, and then, Al was deposited on the electron injection layer to

form a second electrode (cathode) having a thickness of 2,000 Å, thereby completing manufacturing of an organic light-emitting diode.

Example 2

[0254] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound 7 was used instead of Compound 4 in forming the emission layer.

Example 3

[0255] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound 8 was used instead of Compound 4 in forming the emission layer.

Example 4

[0256] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound 13 was used instead of Compound 4 in forming the emission layer.

Example 5

[0257] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound 15 was used instead of Compound 4 in forming the emission layer.

Example 6

[0258] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound 32 was used instead of Compound 4 in forming the emission layer.

Example 7

[0259] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound 35 was used instead of Compound 4 in forming the emission layer.

Example 8

[0260] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound 44 was used instead of Compound 4 in forming the emission layer.

Example 9

[0261] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound 69 was used instead of Compound 4 in forming the emission layer.

Example 10

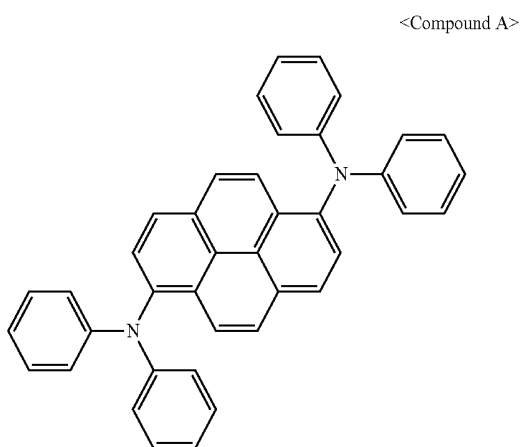
[0262] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound 87 was used instead of Compound 4 in forming the emission layer.

Example 11

[0263] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound 90 was used instead of Compound 4 in forming the emission layer.

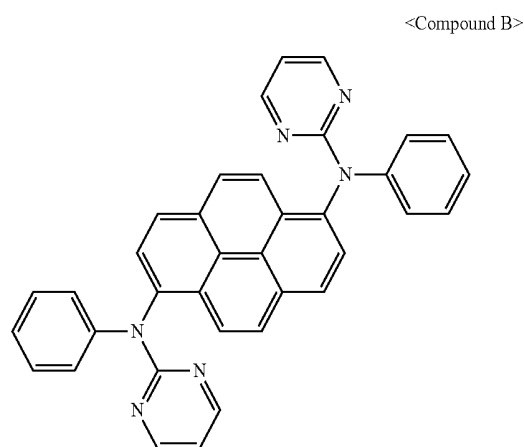
Comparative Example A

[0264] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound A was used instead of Compound 4 in forming the emission layer.



Comparative Example B

[0265] An organic light-emitting diode was manufactured in the same manner as in Example 1, except that as a dopant, Compound B was used instead of Compound 4 in forming the emission layer.



Evaluation Example 2

[0266] The driving voltage, current density, brightness, and a Color coordinate of the organic light-emitting diodes manufactured according to Examples 1 to 11 and a Comparative Examples A and B were evaluated by using a PR650Spectroscan Source Measurement Unit. (product of PhotoResearch Company). Results thereof are shown in Table 2, below. A IS90 lifespan refers to a time required when brightness reduces to 90% of its initial brightness.

TABLE 2

	Emission layer host	Emission layer dopant	Brightness (cd/m ²)	Current density (cd/A)	Driving voltage (V)	Color coordinate (CIE)
Example 1	ADN	Compound 4	700	12	4.2	(0.14, 0.091)
Example 2	ADN	Compound 7	700	13	4.1	(0.14, 0.099)
Example 3	ADN	Compound 8	700	12	4.3	(0.14, 0.085)
Example 4	ADN	Compound 13	700	13	4.3	(0.14, 0.109)
Example 5	ADN	Compound 15	700	13	4.2	(0.14, 0.085)
Example 6	ADN	Compound 32	700	14	4.4	(0.14, 0.095)
Example 7	ADN	Compound 35	700	12	4.3	(0.14, 0.108)
Example 8	ADN	Compound 44	700	13	4.2	(0.14, 0.099)
Example 9	ADN	Compound 69	700	12	4.3	(0.14, 0.101)
Example 10	ADN	Compound 87	700	14	4.2	(0.14, 0.102)
Example 11	ADN	Compound 90	700	12	4.3	(0.14, 0.107)
Comparative Example A	ADN	Compound A	700	16	4.6	(0.16, 0.20)
Comparative Example B	ADN	Compound B	700	44	6.0	(0.16, 0.10)

[0267] From Table 1, it may be seen that the organic light-emitting diodes of Examples 1 to 11 exhibited lower driving voltages, higher current density, and better color purity characteristics than the organic light-emitting diodes of Comparative Examples A and B.

[0268] An organic light-emitting diode and an organic light-emitting apparatus each including the pyrene-based compound according to an embodiment may have a low driving voltage, a high current density, and excellent color purity characteristics, and a long lifespan.

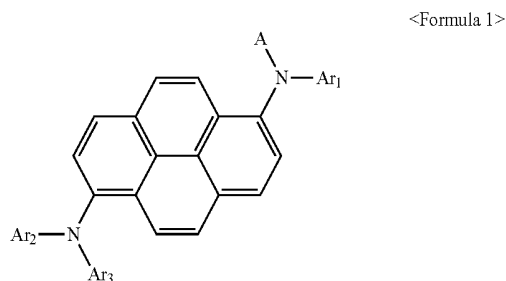
[0269] By way of summation and review, an OLED may have a structure including a substrate, and an anode, a hole transport layer, an emission layer, an electron transport layer, and a cathode which are sequentially stacked on the substrate. The hole transport layer, the emission layer, and the electron transport layer may be organic thin films formed of organic compounds.

[0270] An operating principle of an OLED having the above-described structure may be as follows. Holes injected from the anode may move to the emission layer via the hole transport layer, and electrons injected from the cathode may move to the emission layer via the electron transport layer. The holes and electrons (carriers) may recombine in the organic emission layer to generate excitons. When the excitons drop from an excited state to a ground state, light is emitted.

[0271] Example embodiments have been disclosed herein, and although specific terms are employed, they are used and are to be interpreted in a generic and descriptive sense only and not for purpose of limitation. In some instances, as would be apparent to one of ordinary skill in the art as of the filing of the present application, features, characteristics, and/or elements described in connection with a particular embodiment may be used singly or in combination with features, characteristics, and/or elements described in connection with other embodiments unless otherwise specifically indicated. Accordingly, it will be understood by those of skill in the art that various changes in form and details may be made without departing from the spirit and scope of the present invention as set forth in the following claims.

What is claimed is:

1. A pyrene-based compound represented by Formula 1, below:



wherein, in Formula 1,

A is a substituted or unsubstituted C₂-C₃₀ heteroaromatic group including at least one nitrogen (N); and

Ar₁, Ar₂, and Ar₃ are each independently a substituted or unsubstituted phenyl group or a substituted or unsubstituted naphthyl group.

2. The compound of claim 1, wherein A is selected from a substituted or unsubstituted pyridinyl group, a substituted or

unsubstituted pyrimidinyl group, a substituted or unsubstituted triazinyl group, a substituted or unsubstituted pyrazinyl group, a substituted or unsubstituted pyridazinyl group, a substituted or unsubstituted isindolyl group, a substituted or unsubstituted indolyl group, a substituted or unsubstituted indazolyl group, a substituted or unsubstituted isoquinolinyl group, a substituted or unsubstituted quinolinyl group, a substituted or unsubstituted phthalazinyl group, a substituted or unsubstituted quinoxalinyl group, a substituted or unsubstituted quinazolinyl group, a substituted or unsubstituted cinnolinyl group, and a substituted or unsubstituted 1,10-phenanthrolyl group.

3. The organic light-emitting diode of claim 1, wherein A is selected from:

a pyridinyl group, a pyrimidinyl group, a triazinyl group, a pyrazinyl group, a pyridazinyl group, an isindolyl group, an indolyl group, an indazolyl group, an isoquinolinyl group, a quinolinyl group, a phthalazinyl group, a quinoxalinyl group, a cinnolinyl group, and a 1,10-phenanthrolyl group; or

a pyridinyl group, a pyrimidinyl group, a triazinyl group, a pyrazinyl group, a pyridazinyl group, an isindolyl group, an indolyl group, an indazolyl group, an isoquinolinyl group, a quinolinyl group, a phthalazinyl group, a quinoxalinyl group, a cinnolinyl group, and a 1,10-phenanthrolyl group, each substituted with at least one of:

a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group,

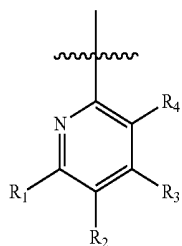
a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, and a C₁-C₆₀ alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof,

a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, and a C₂-C₆₀ heteroaryl group,

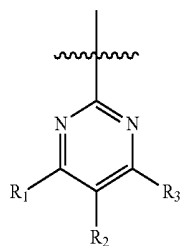
a C₃-C₁₀ cycloalkyl group, a C₃-C₁₀ heterocycloalkyl group, a C₃-C₁₀ cycloalkenyl group, a C₃-C₁₀ heterocycloalkenyl group, a C₆-C₆₀ aryl group, a C₆-C₆₀ aryloxy group, a C₆-C₆₀ arylthio group, and a C₂-C₆₀ heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₆₀ alkyl group, a C₂-C₆₀ alkenyl group, a C₂-C₆₀ alkynyl group, a C₁-C₆₀ alkoxy group, a phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group, or

—Si(Q₁₁)(Q₁₂)(Q₁₃), in which Q₁₁ to Q₁₃ are each independently a C₁-C₆₀ alkyl group, a C₁-C₆₀ alkoxy group, a C₆-C₆₀ aryl group, or a C₂-C₆₀ heteroaryl group.

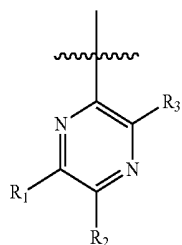
4. The compound of claim 1, wherein A is a group represented by one of Formulae 2A to 2K below:



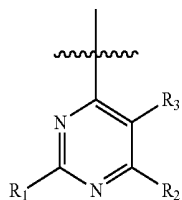
Formula 2A



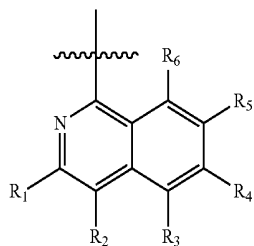
Formula 2B



Formula 2C

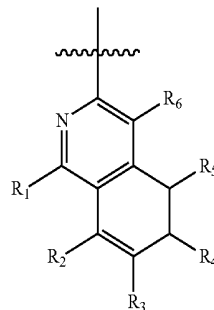


Formula 2D

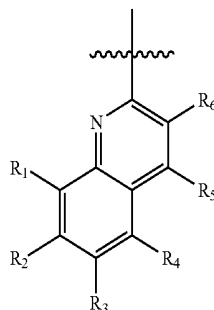


Formula 2E

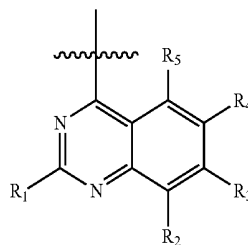
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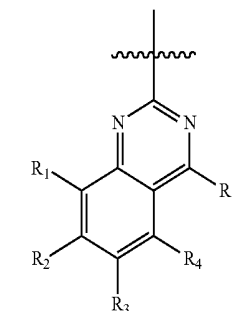
Formula 2F



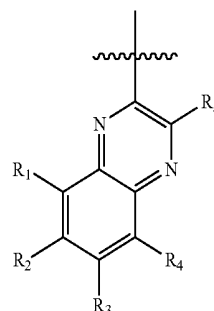
Formula 2G



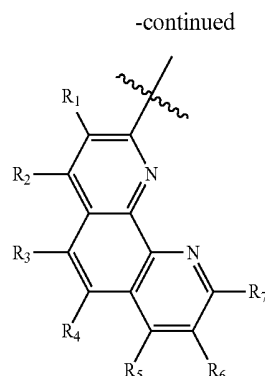
Formula 2H



Formula 2I



Formula 2J



Formula 2K

wherein, in Formulae 2A to 2K, R_1 to R_7 are each independently selected from:

- a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group;
- a C_1 - C_{60} alkyl group, a C_2 - C_{60} alkenyl group, a C_2 - C_{60} alkynyl group, and a C_1 - C_{60} alkoxy group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof;
- a C_3 - C_{10} cycloalkyl group, a C_3 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_3 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, and a C_2 - C_{60} heteroaryl group;
- a C_3 - C_{10} cycloalkyl group, a C_3 - C_{10} heterocycloalkyl group, a C_3 - C_{10} cycloalkenyl group, a C_3 - C_{10} heterocycloalkenyl group, a C_6 - C_{60} aryl group, a C_6 - C_{60} aryloxy group, a C_6 - C_{60} arylthio group, and a C_2 - C_{60} heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic 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, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group; and
- Si(Q_{11})(Q_{12})(Q_{13}), in which Q_{11} to Q_{13} are each independently a C_1 - C_{10} alkyl group, a C_6 - C_{60} aryl group, or a C_2 - C_{60} heteroaryl group.

5. The pyrene-based compound of claim 4, wherein R_1 to R_7 are each independently selected from:

- a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a C_1 - C_{10} alkyl group;

- a C_1 - C_{60} alkyl group substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof;

- a C_6 - C_{16} aryl group, and a C_2 - C_{16} heteroaryl group;

- a C_6 - C_{16} aryl group or a C_2 - C_{16} heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C_1 - C_{10} alkyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenylcarbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group; and

- Si(Q_{11})(Q_{12})(Q_{13}), in which Q_{11} to Q_{13} are each independently a C_1 - C_{10} alkyl group, a C_6 - C_{16} aryl group, or a C_2 - C_{16} heteroaryl group.

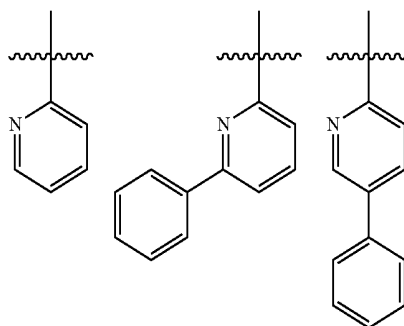
6. The pyrene-based compound of claim 4, wherein R_1 to R_7 are each independently selected from:

- a hydrogen atom, a deuterium atom, F, Cl, Br, I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid 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, and a hexyl group;

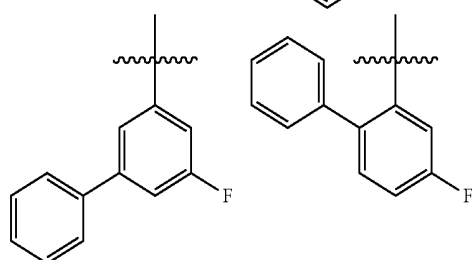
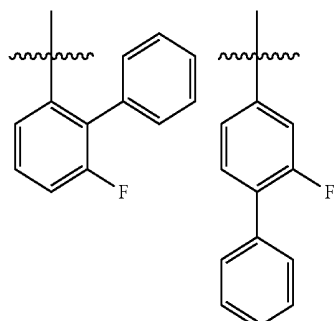
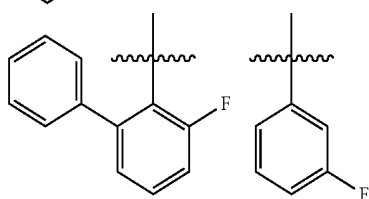
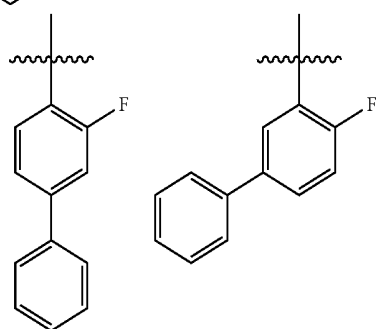
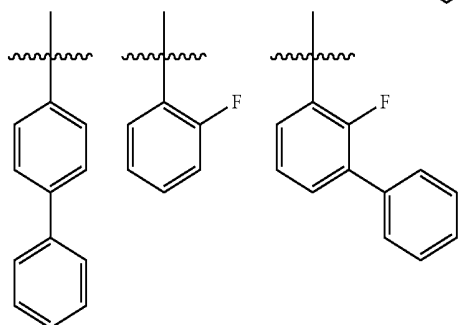
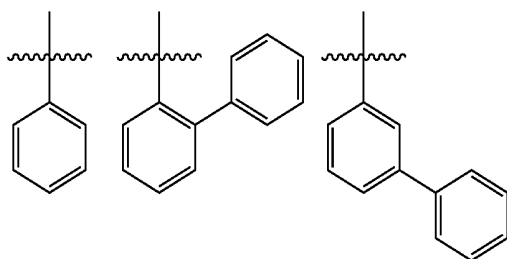
- a phenyl group, a naphthyl group, and an anthracenyl group; and

- a phenyl group, a naphthyl group, and an anthracenyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a methyl group, an ethyl group, a propyl group, a butyl group, a pentyl group, a hexyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, and a phenylcarbazolyl group.

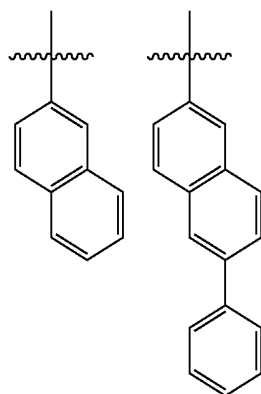
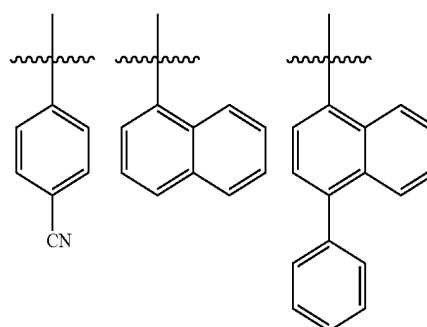
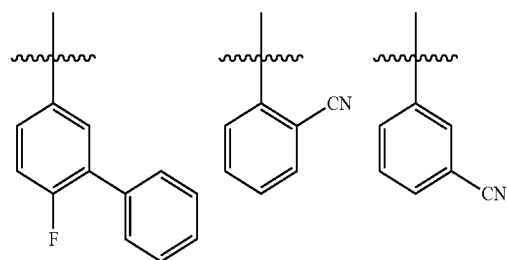
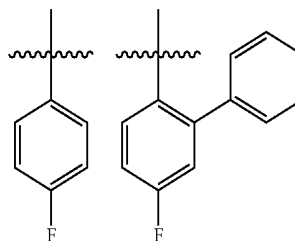
7. The pyrene-based compound of claim 1, wherein A has a structure selected from one of the following structures:



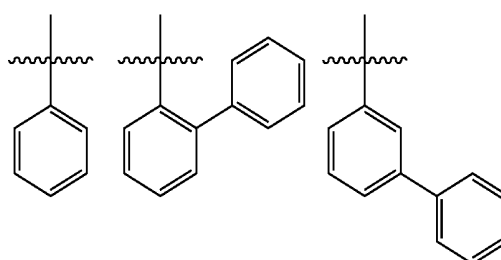
10. The pyrene-based compound of claim 1, wherein Ar₁ is selected from one of the following structures:

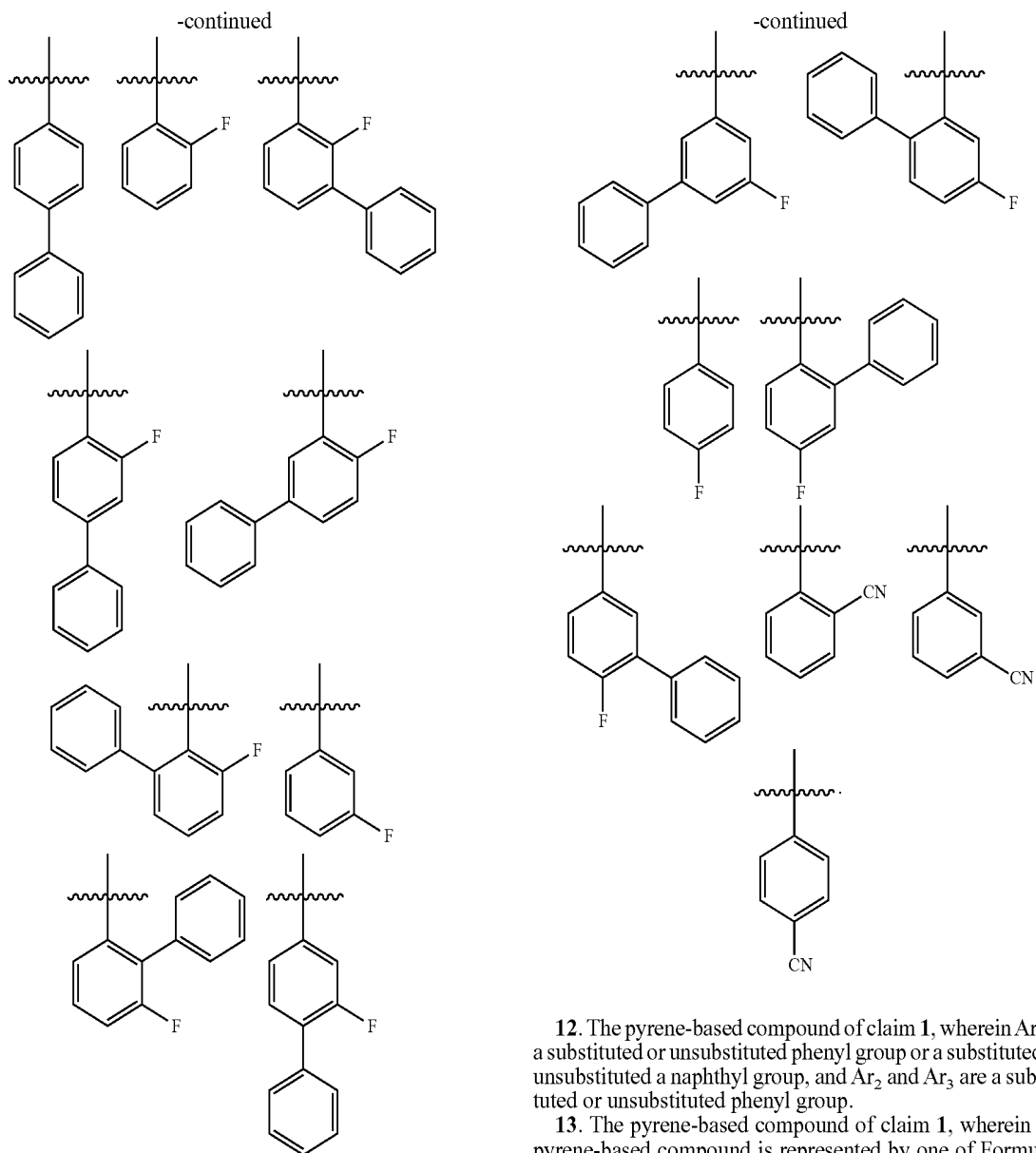


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11. The pyrene-based compound of claim 1, wherein Ar₂ and Ar₃ are each independently selected from one of the following structures:

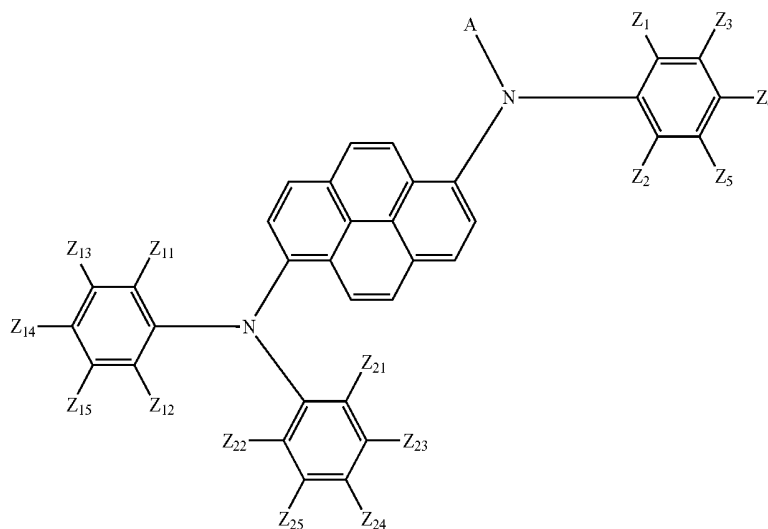




12. The pyrene-based compound of claim 1, wherein Ar_1 is a substituted or unsubstituted phenyl group or a substituted or unsubstituted naphthyl group, and Ar_2 and Ar_3 are a substituted or unsubstituted phenyl group.

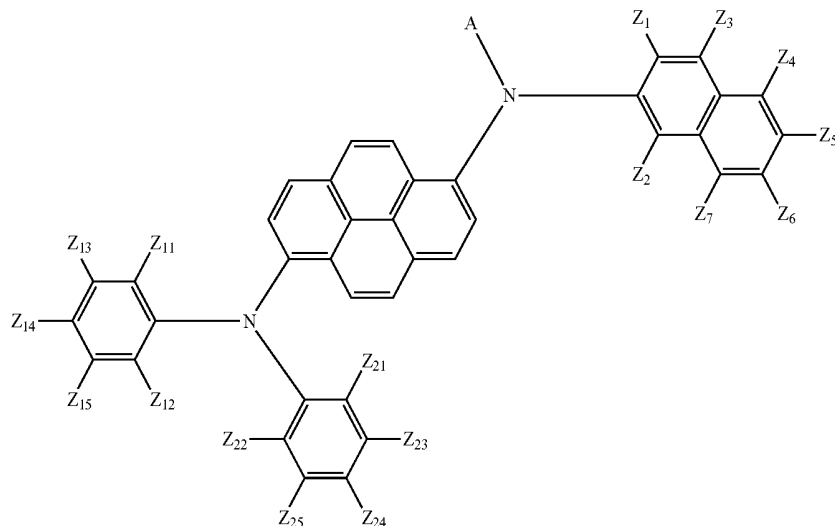
13. The pyrene-based compound of claim 1, wherein the pyrene-based compound is represented by one of Formulae 1A, 1B, or 1C, below:

<Formula 1A>

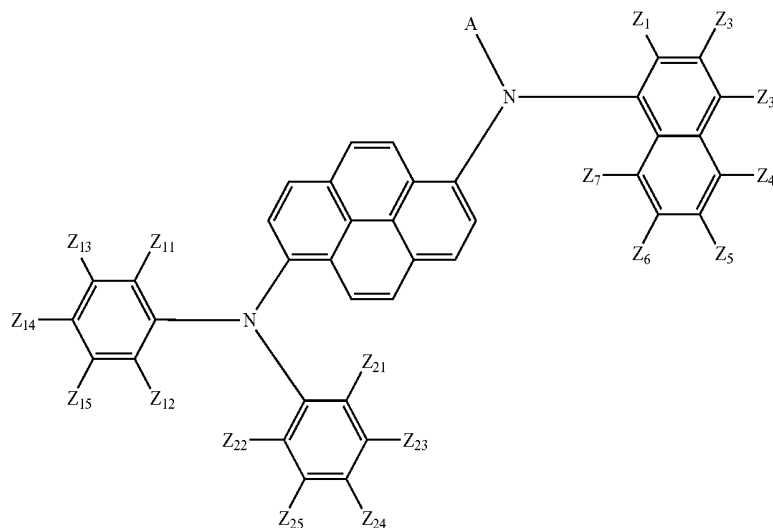


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<Formula 1B>



<Formula 1C>



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

A is a substituted or unsubstituted C_2 - C_{30} heteroaromatic group including at least one nitrogen (N); and

Z_1 to Z_7 , Z_{11} to Z_{15} and Z_{21} to Z_{25} are each independently selected from:

a hydrogen atom, a deuterium atom, F, Cl, Br, I, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid 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, and a hexyl group;

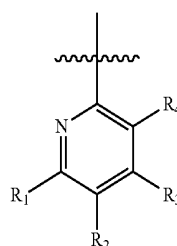
a phenyl group, a naphthyl group, and an anthracenyl group; and

a phenyl group, a naphthyl group, and an anthracenyl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a methyl

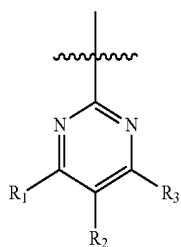
group, an ethyl group, a propyl group, a butyl group, a pentyl group, a hexyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group and a phenylcarbazolyl group.

14. The pyrene-based compound of claim 13, wherein A is represented by one of Formulae 2A to 2K, below:

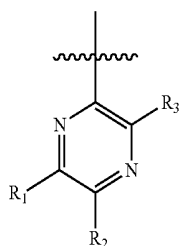
Formula 2A



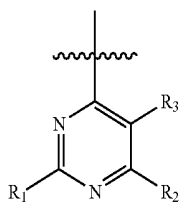
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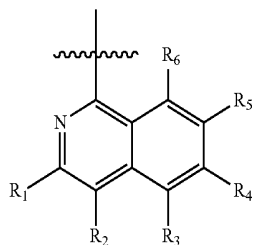
Formula 2B



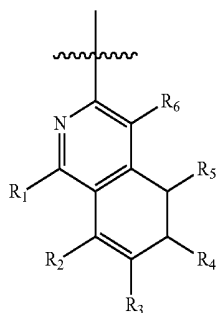
Formula 2C



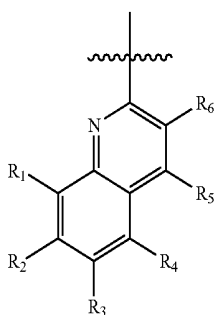
Formula 2D



Formula 2E

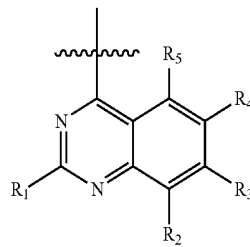


Formula 2F

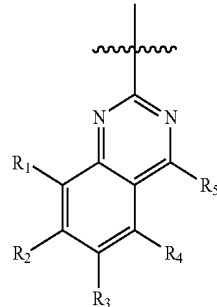


Formula 2G

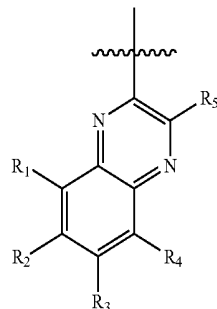
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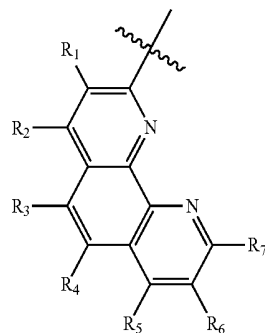
Formula 2H



Formula 2I



Formula 2J



Formula 2K

wherein, in Formulae 2A to 2K, R_1 to R_7 are each independently selected from:

- a hydrogen atom, a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a C_1 - C_{10} alkyl group;
- a C_1 - C_{60} alkyl group substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, and a phosphoric acid or a salt thereof;

a C₆-C₁₆ aryl group and a C₂-C₁₆ heteroaryl group;

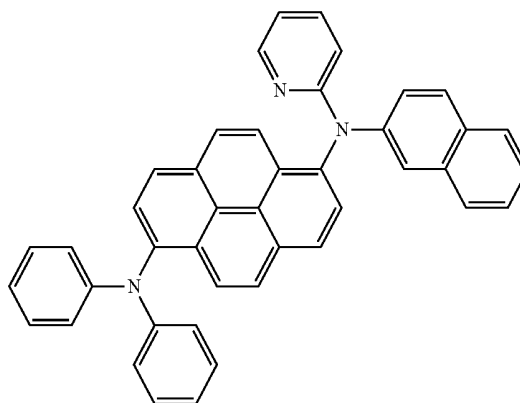
a C₆-C₁₆ aryl group and a C₂-C₁₆ heteroaryl group, each substituted with at least one of a deuterium atom, a halogen atom, a hydroxyl group, a cyano group, a nitro group, an amino group, an amidino group, a hydrazine group, a hydrazone group, a carboxyl group or a salt thereof, a sulfonic acid or a salt thereof, a C₁-C₁₀ alkyl group, phenyl group, a naphthyl group, an anthracenyl group, a fluorenyl group, a dimethylfluorenyl group, a diphenylfluorenyl group, a carbazolyl group, a phenyl-carbazolyl group, a pyridinyl group, a pyrimidinyl group, a pyrazinyl group, a pyridazinyl group, a triazinyl group, a quinolinyl group, and an isoquinolinyl group; and

—Si(Q₁₁)(Q₁₂)(Q₁₃), in which Q₁₁ to Q₁₃ are each independently a C₁-C₁₀ alkyl group, a C₆-C₁₆ aryl group, or a C₂-C₁₆ heteroaryl group.

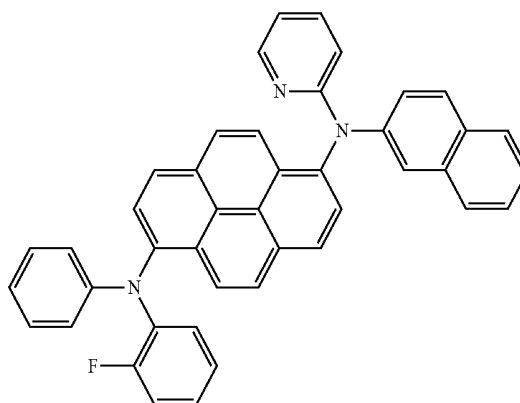
15. The pyrene-based compound of claim **14**, wherein R₁ to R₇, Z₁ to Z₇, Z₁₁ to Z₁₅ and Z₂₁ to Z₂₅ are each independently a hydrogen atom, a F atom, a cyano group, or a phenyl group.

16. The pyrene-based compound of claim **1**, wherein the pyrene-based compound is one of Compounds 1 to 96, below:

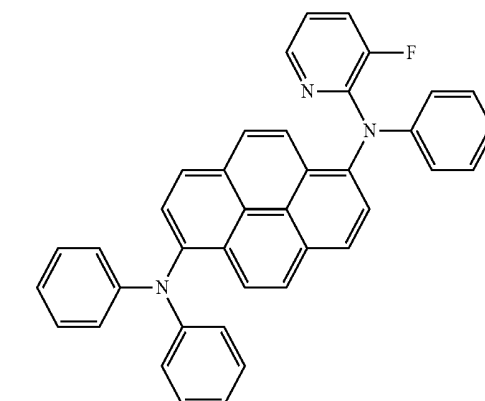
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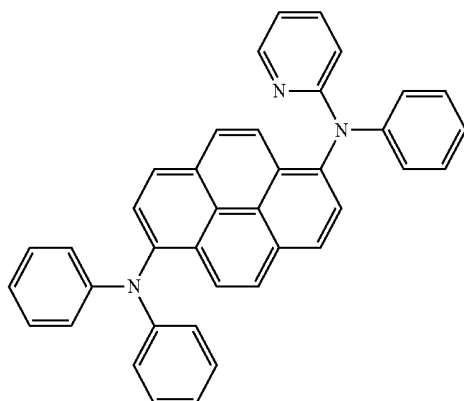
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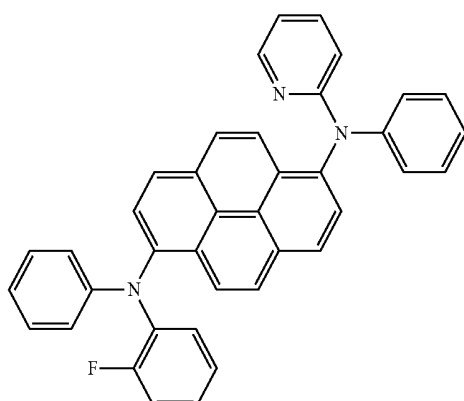
4



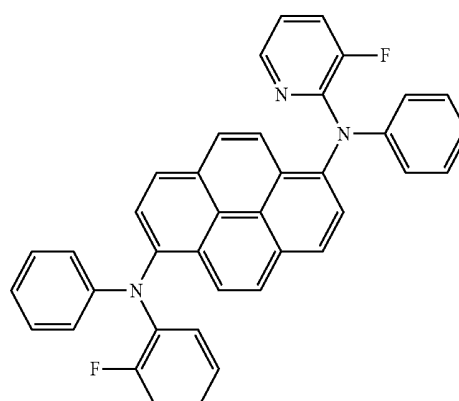
5



1

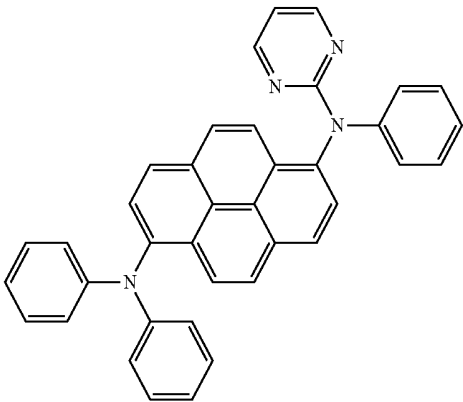


2

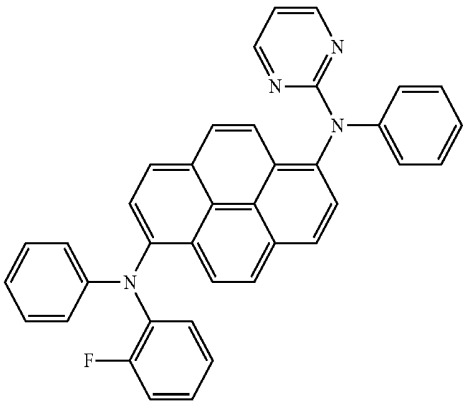


6

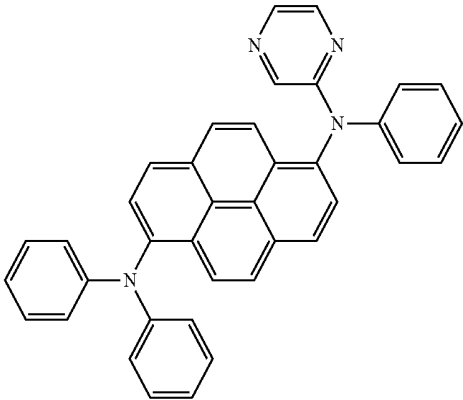
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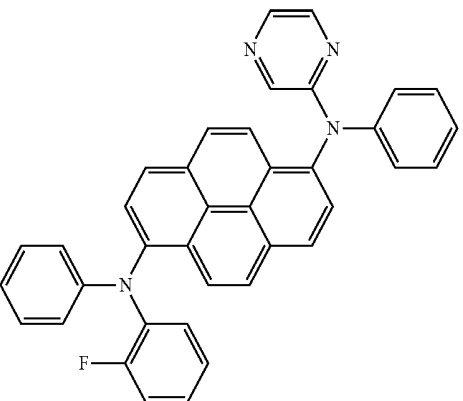
7



8

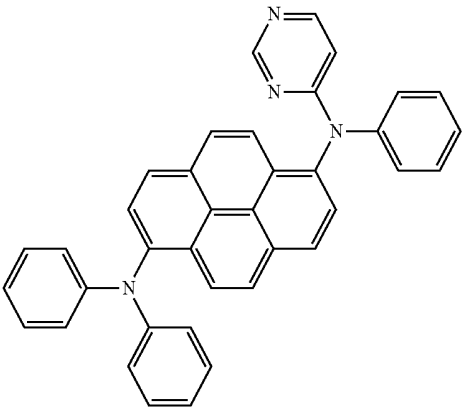


9

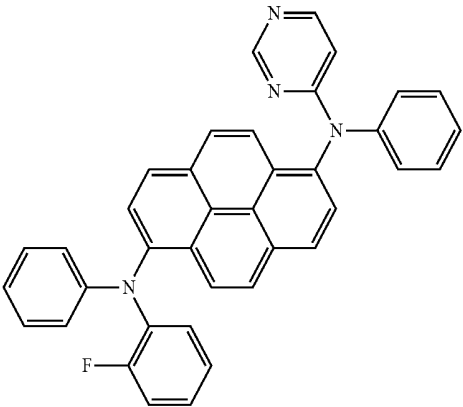


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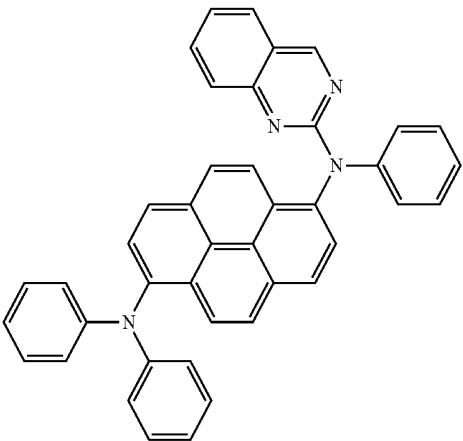
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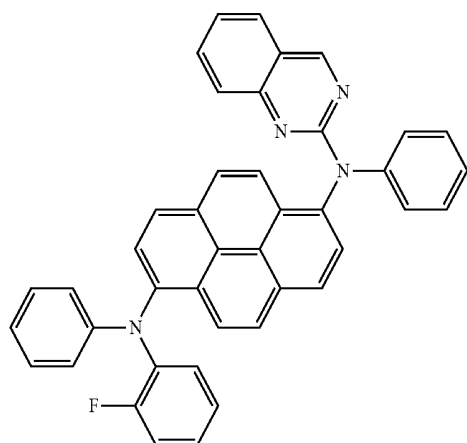


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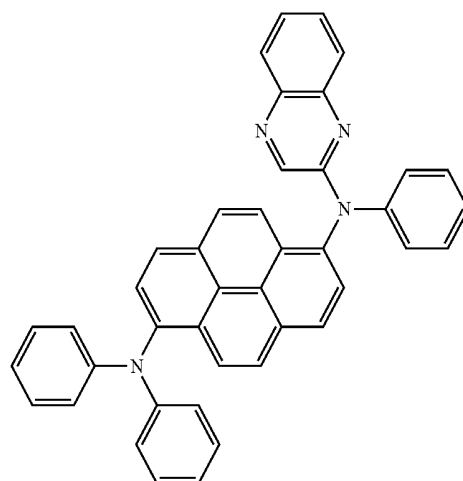
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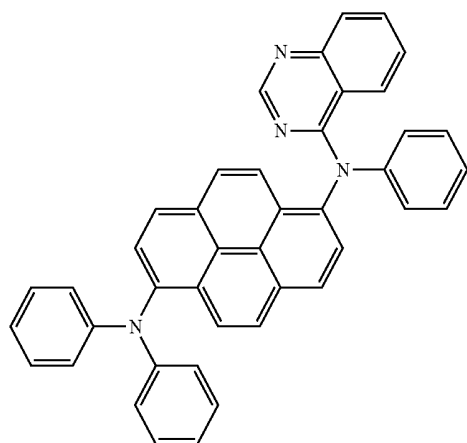


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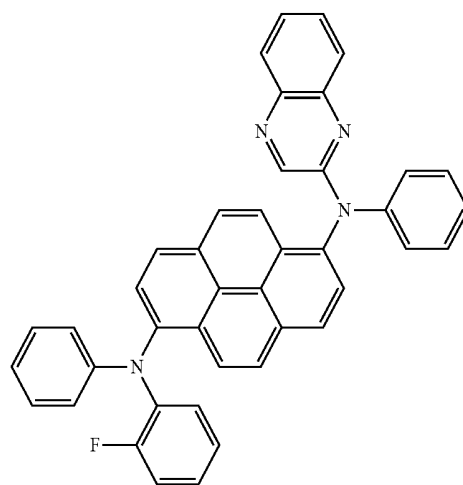
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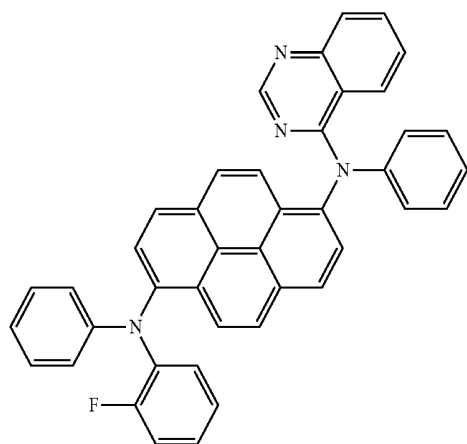
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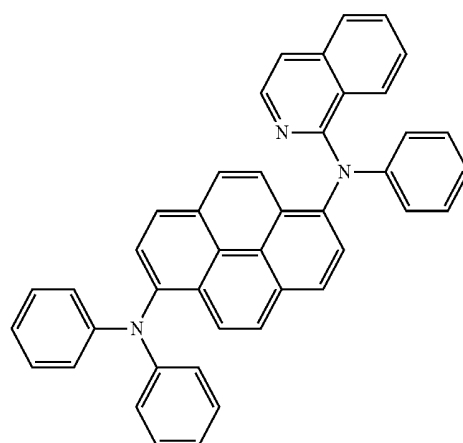
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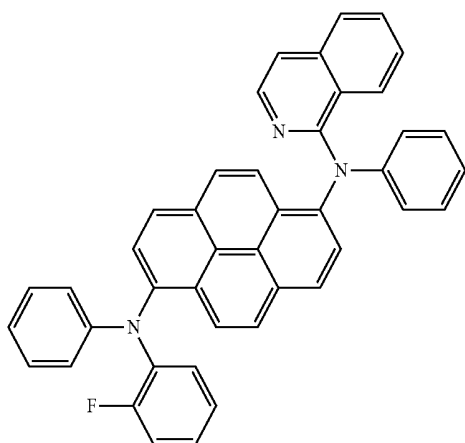


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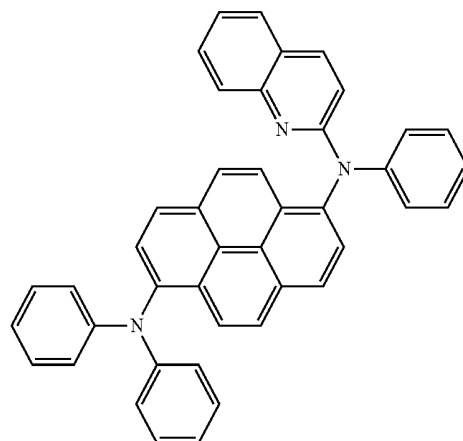
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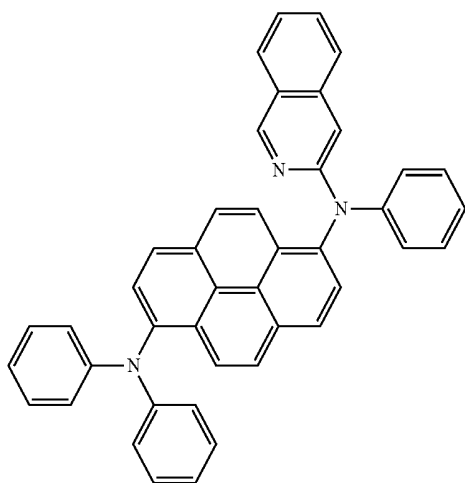


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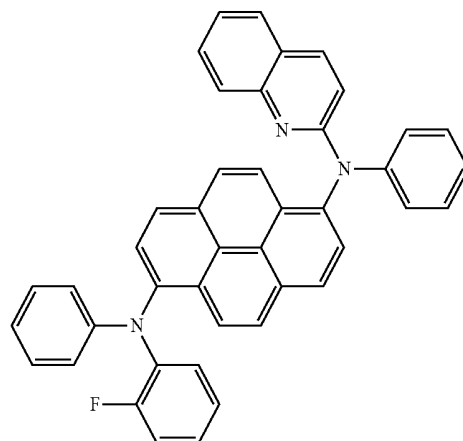
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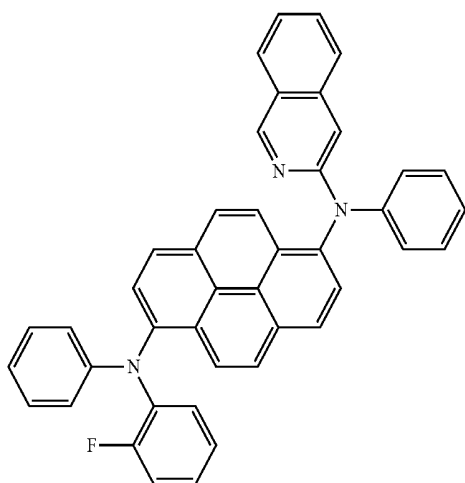
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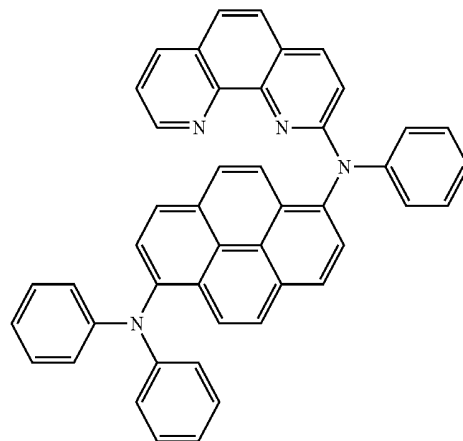
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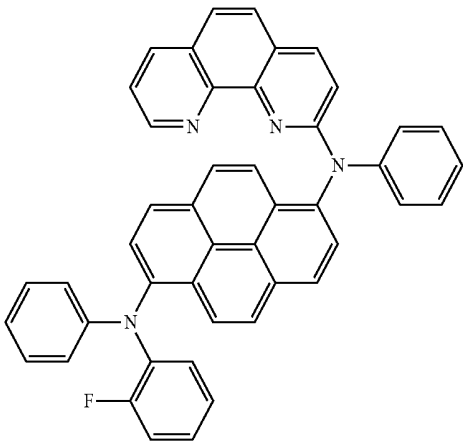


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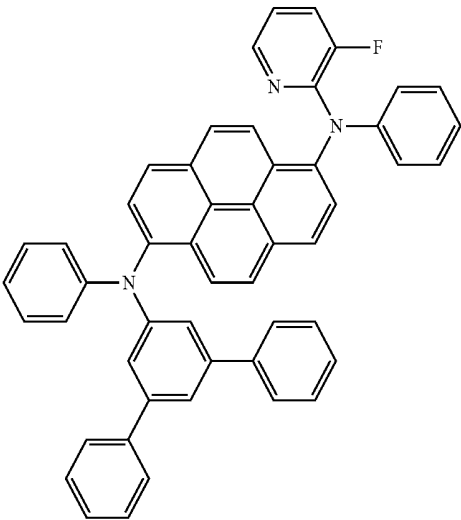
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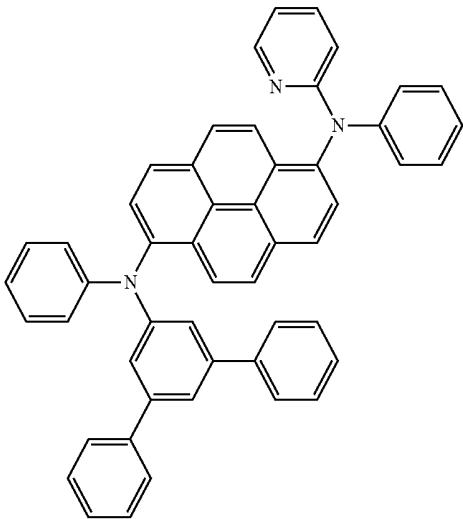


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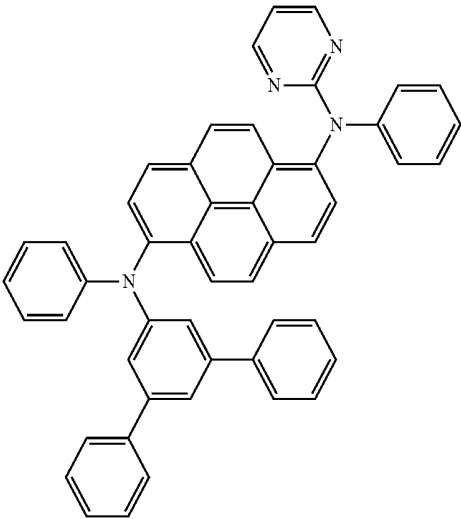
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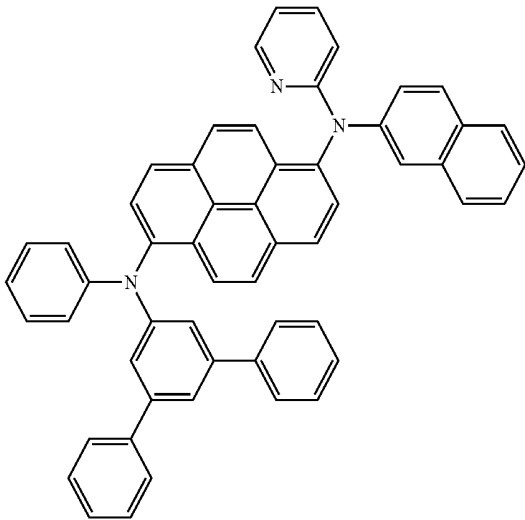
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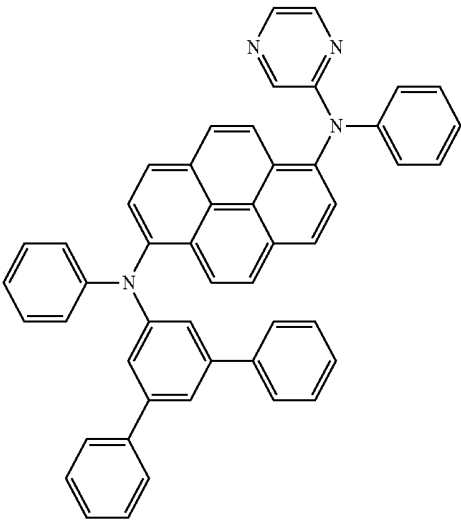
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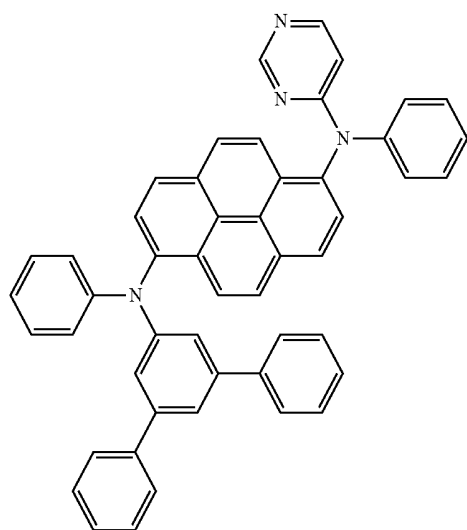
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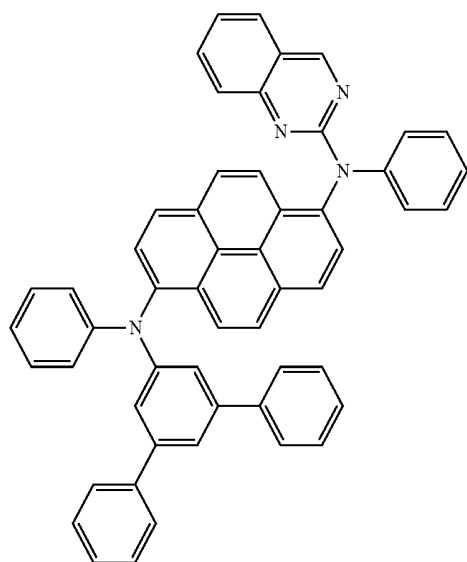
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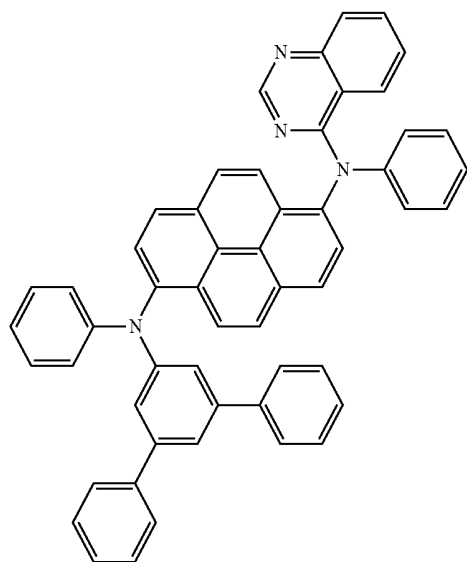
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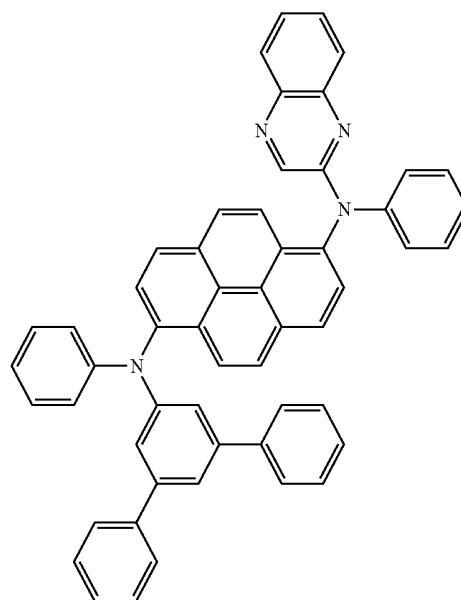


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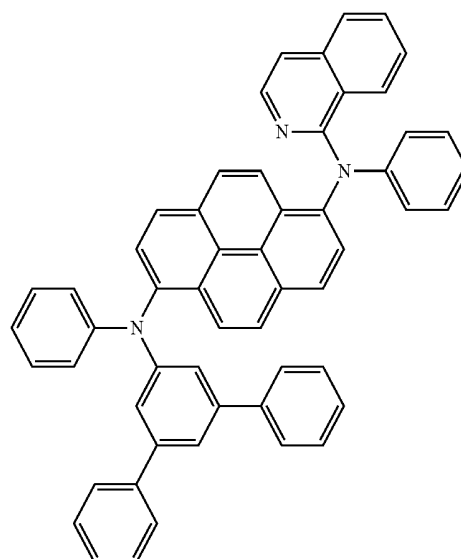


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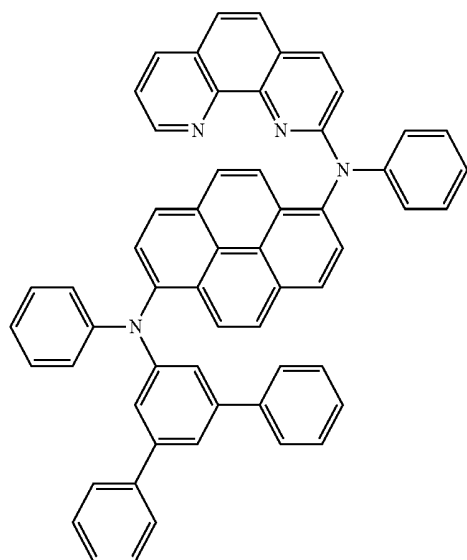
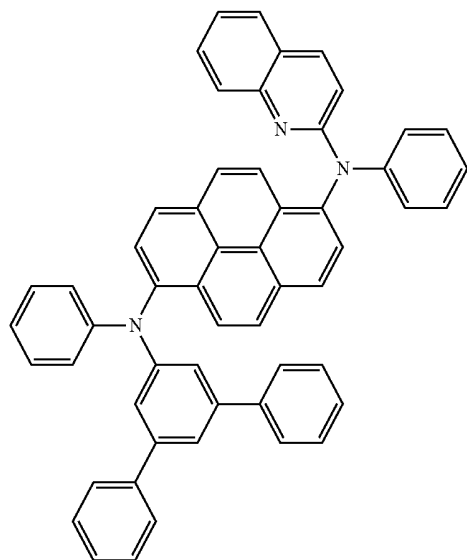
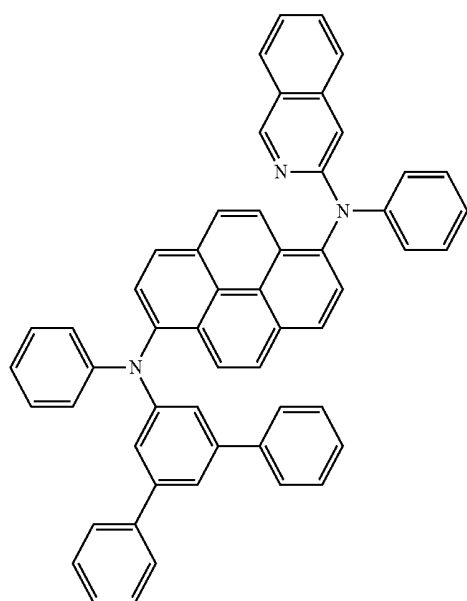
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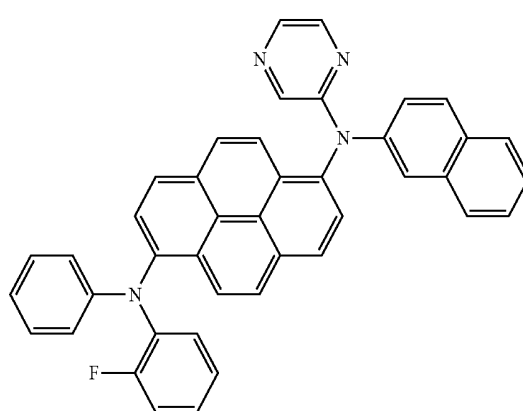
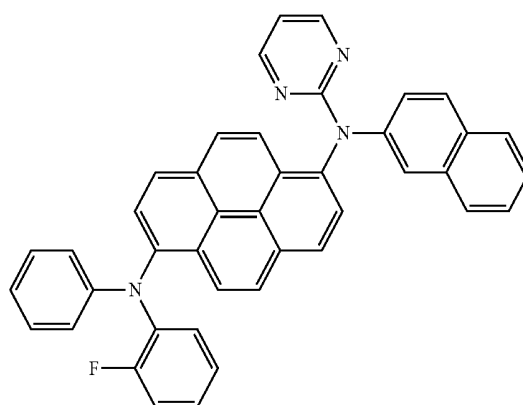
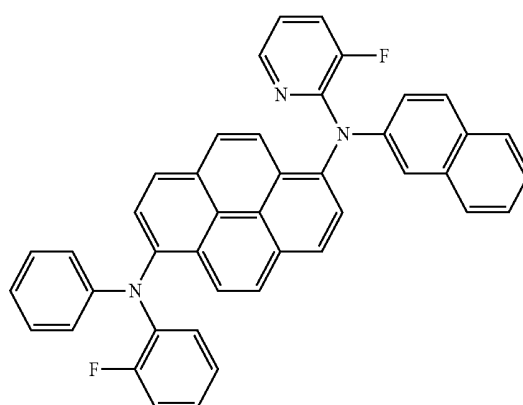
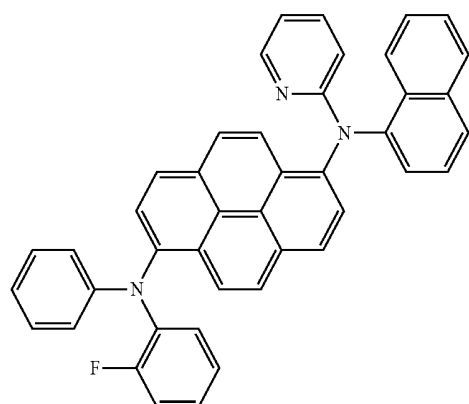
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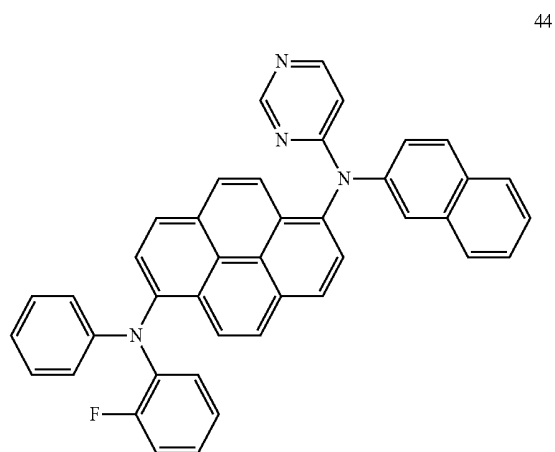
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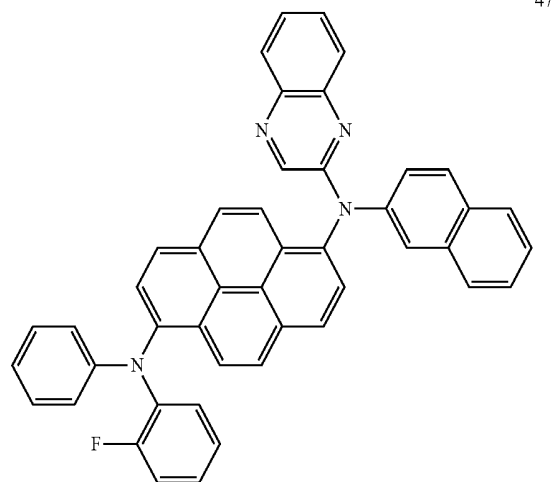
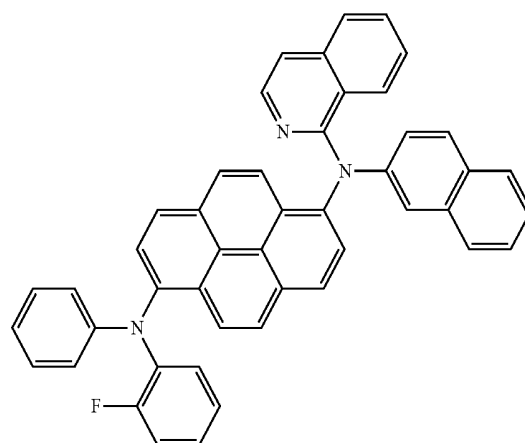
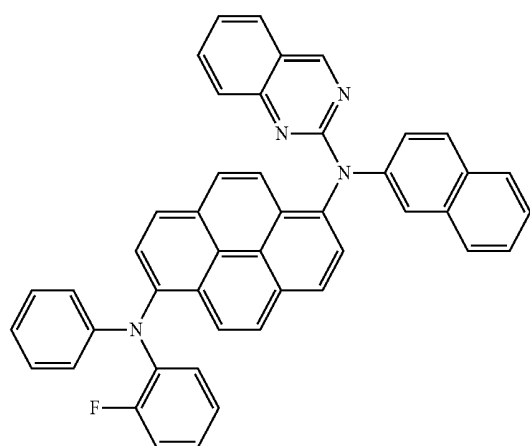
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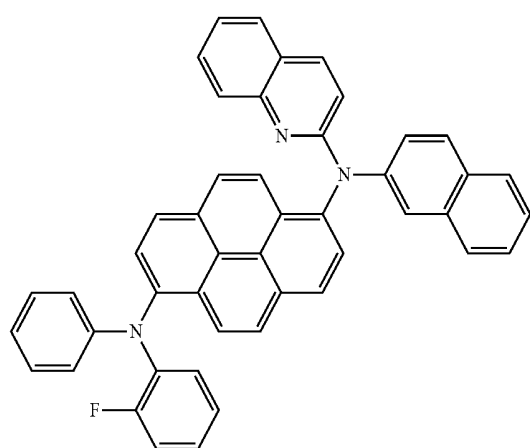
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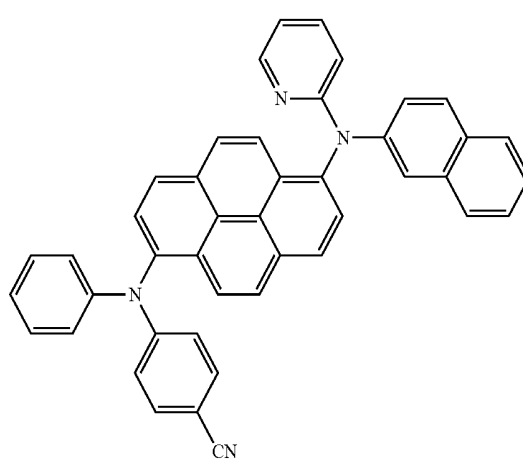
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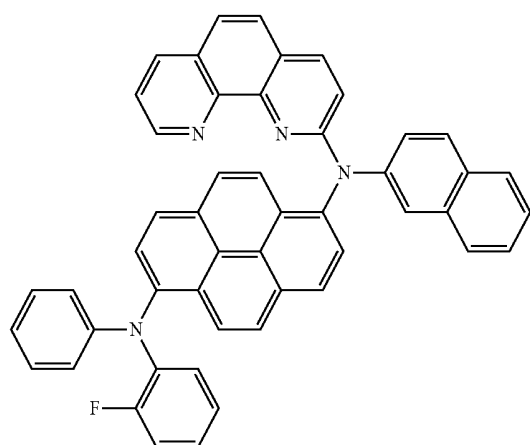
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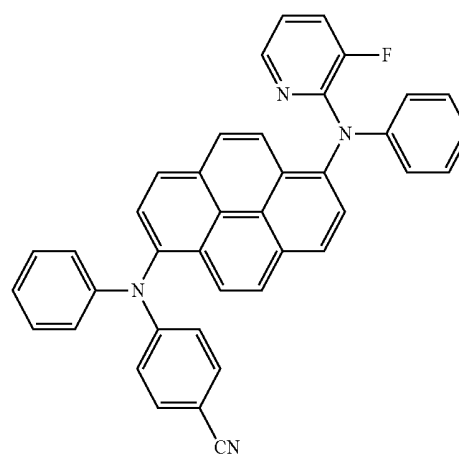
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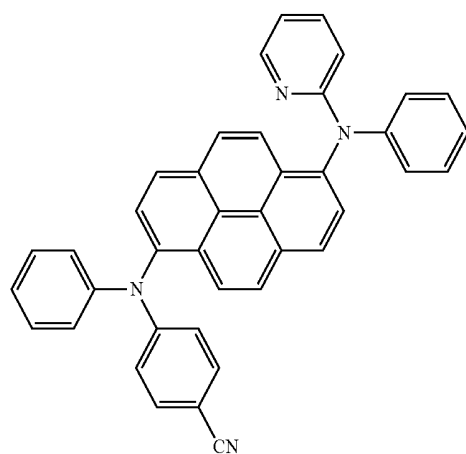
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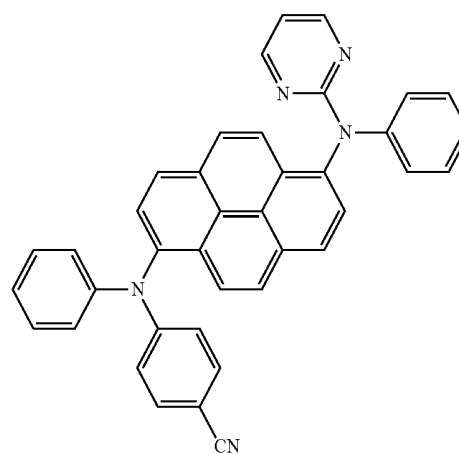
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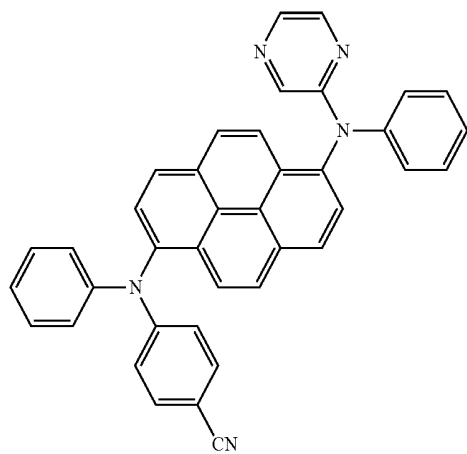
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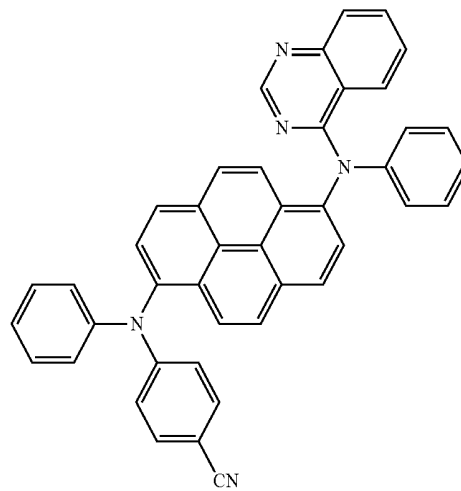


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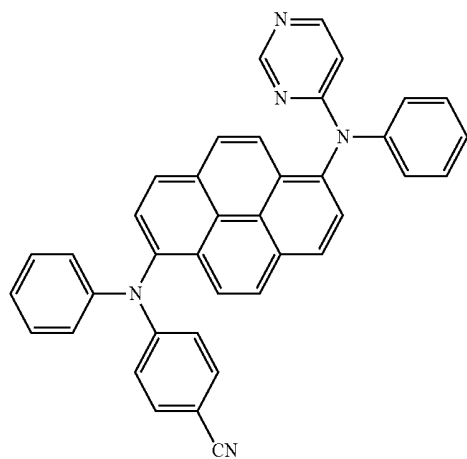


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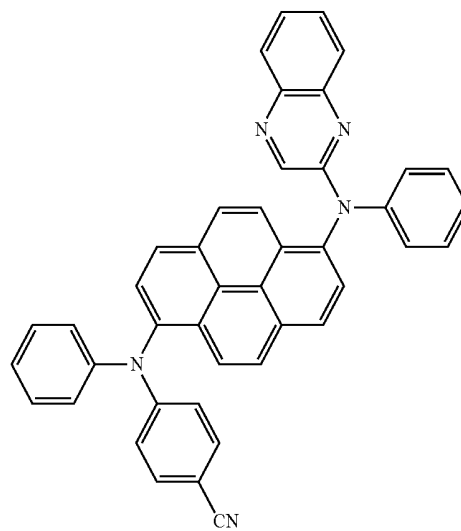
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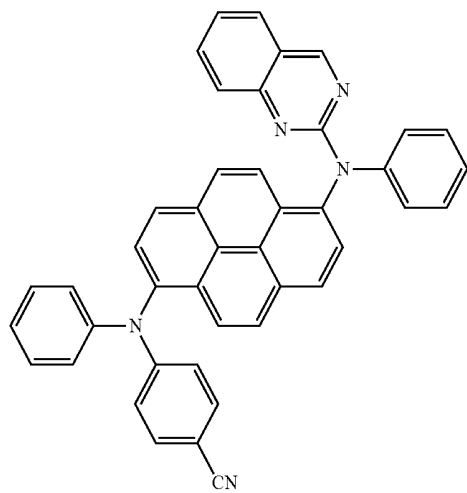
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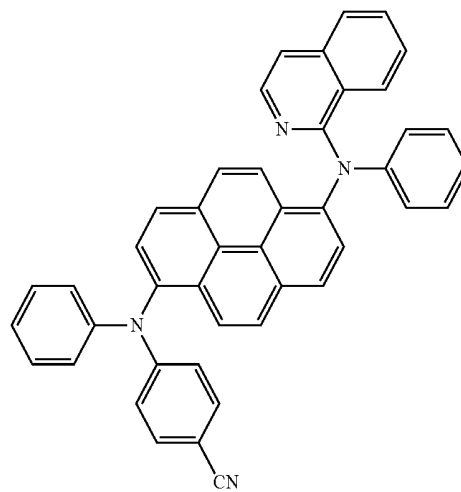
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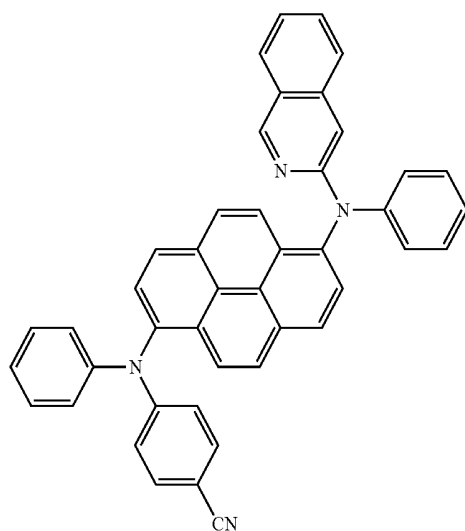


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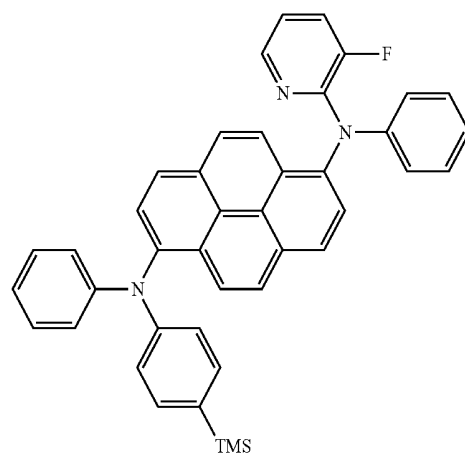
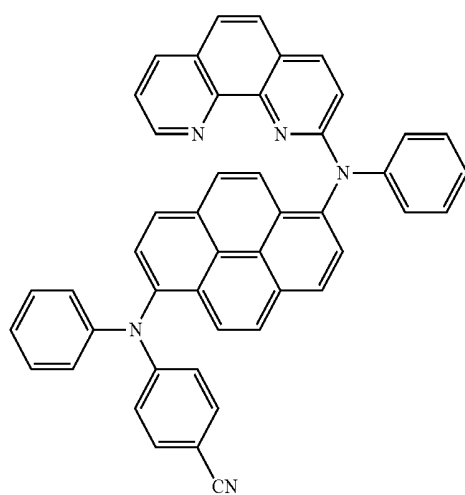
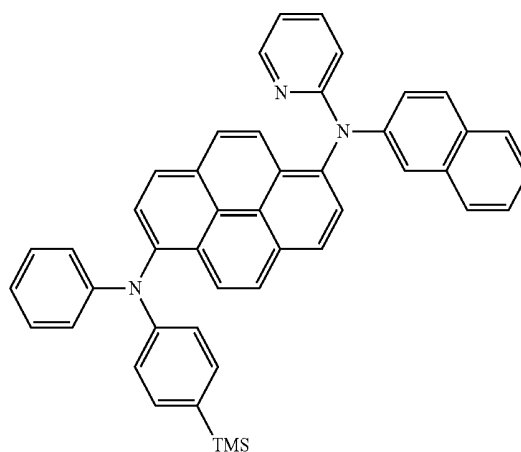
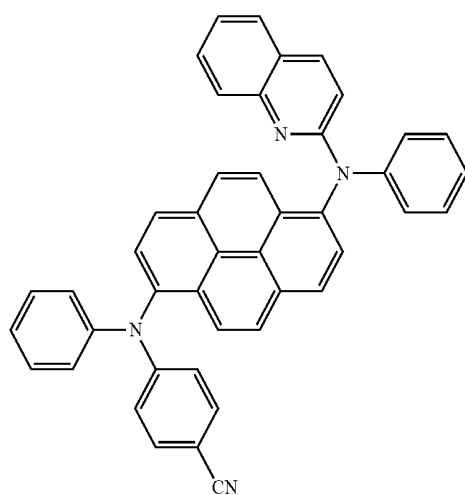
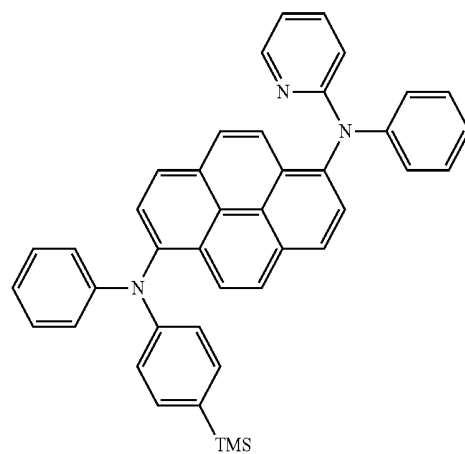


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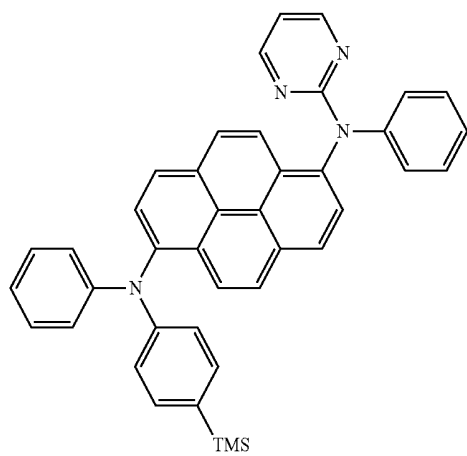
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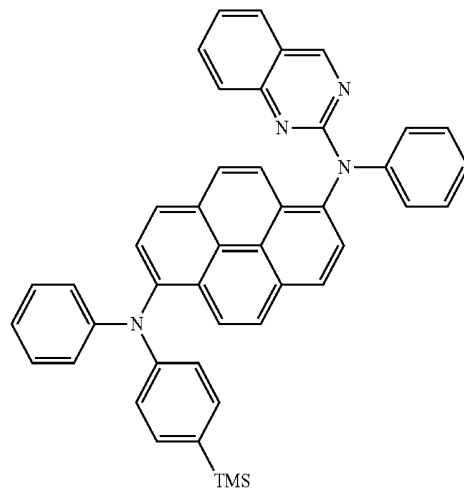
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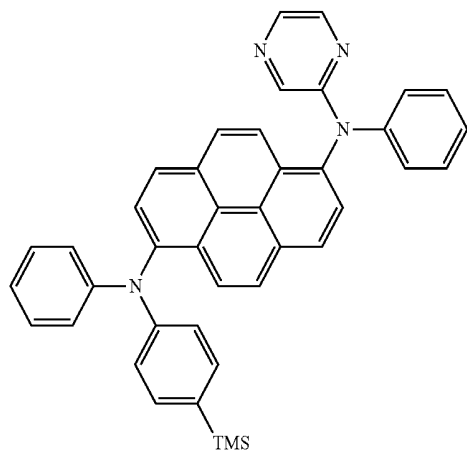
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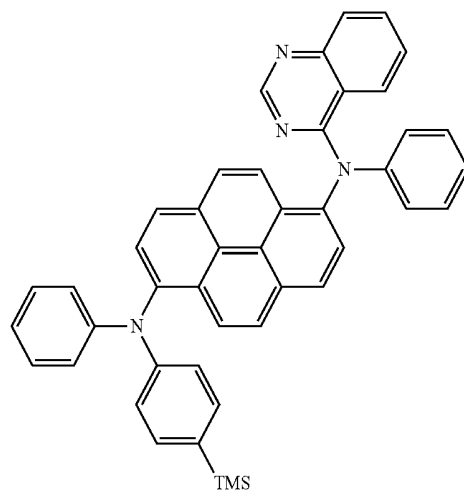
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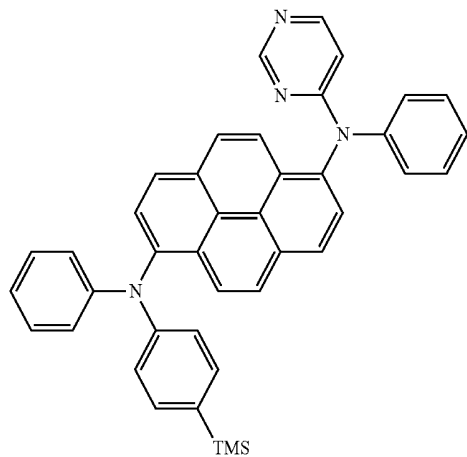
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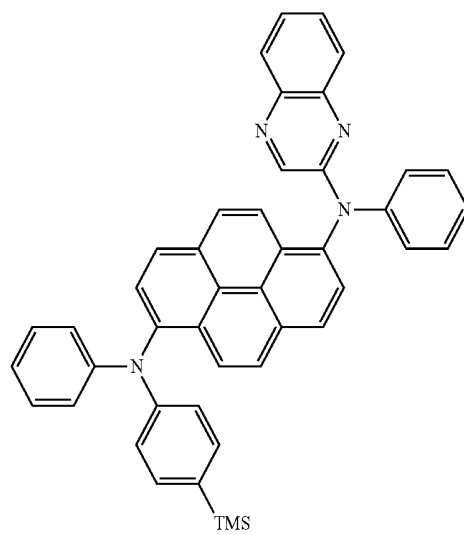
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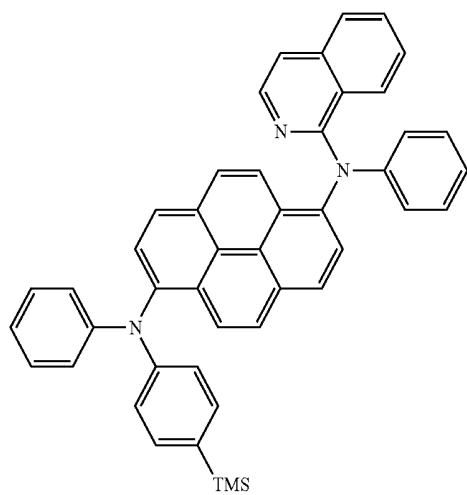
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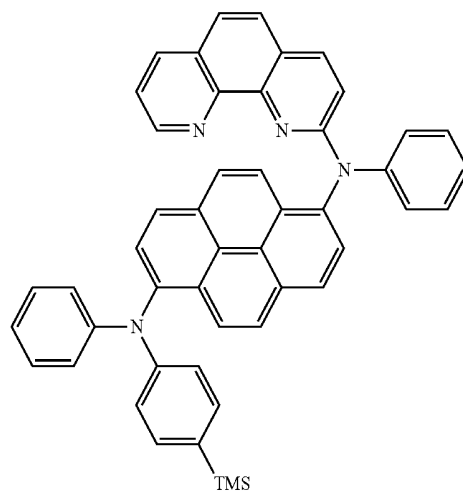


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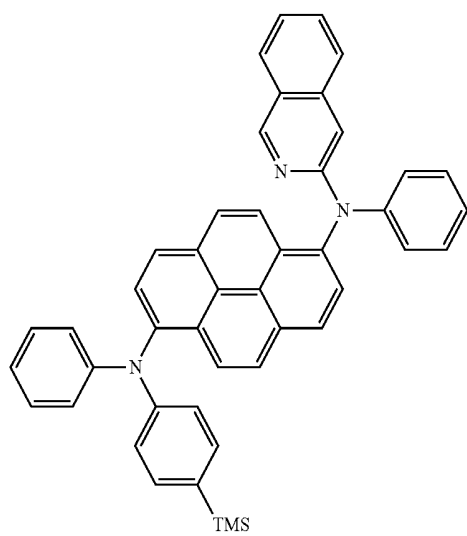


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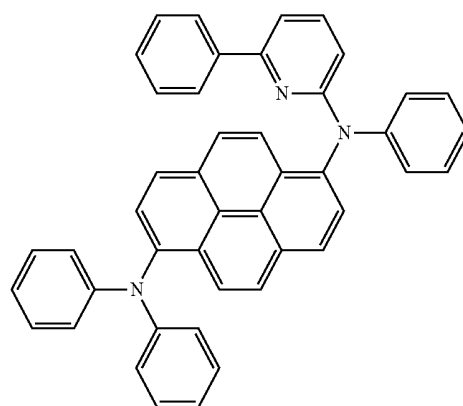
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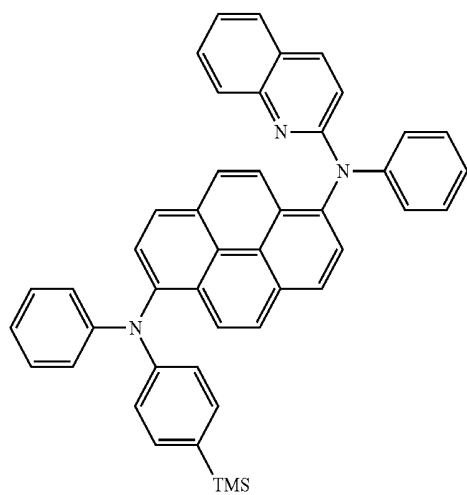
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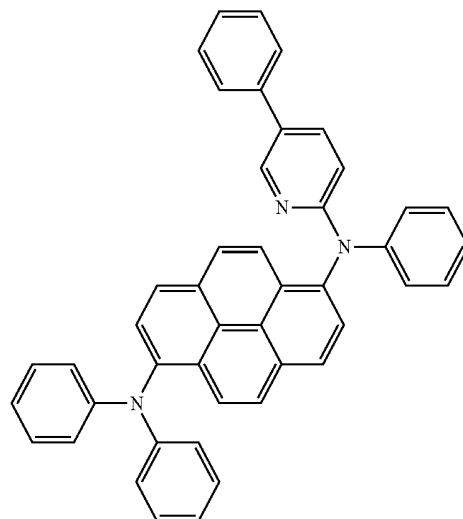
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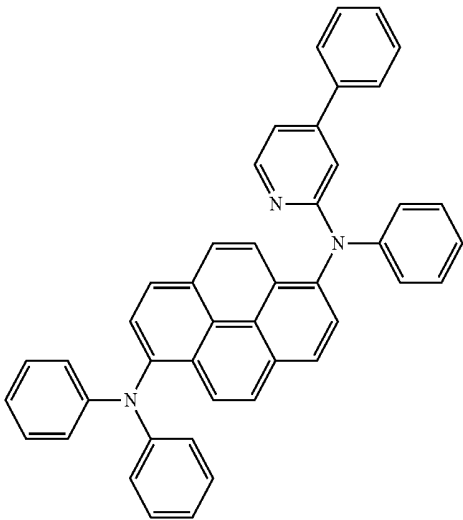


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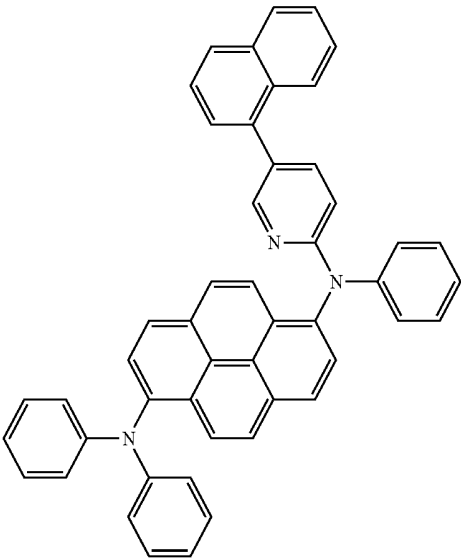
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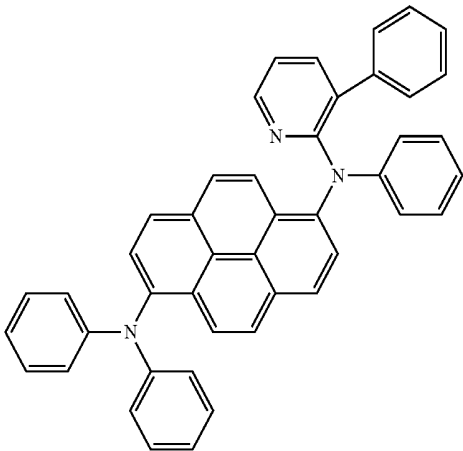


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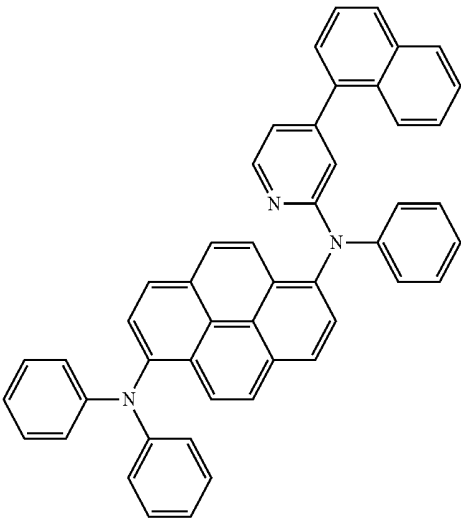
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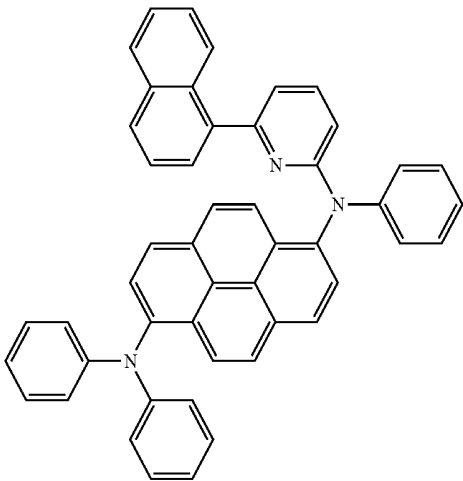
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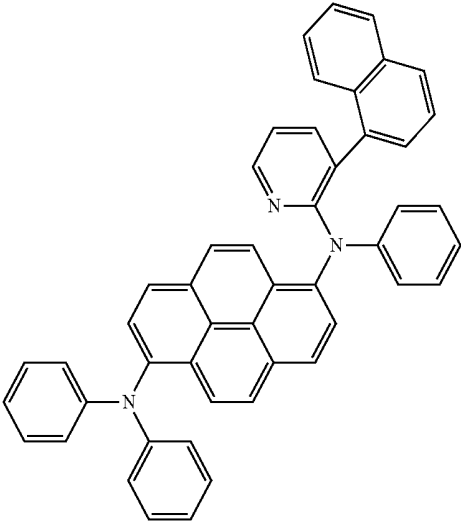
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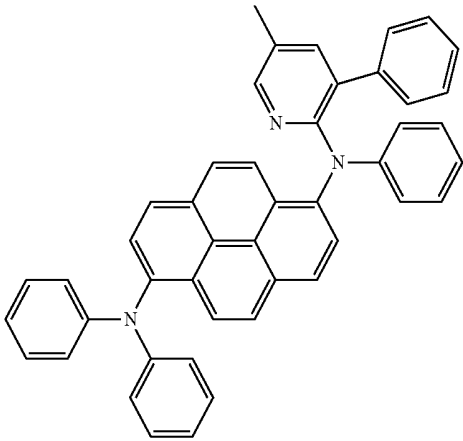


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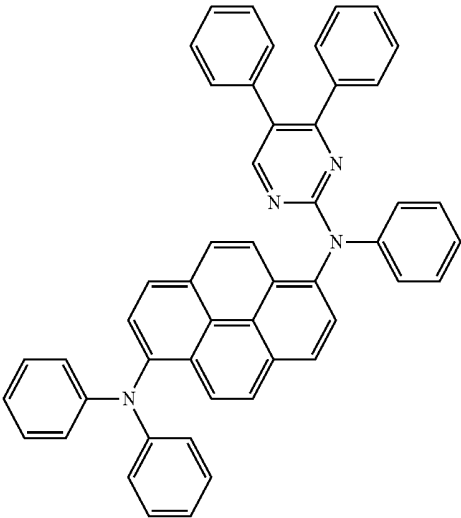
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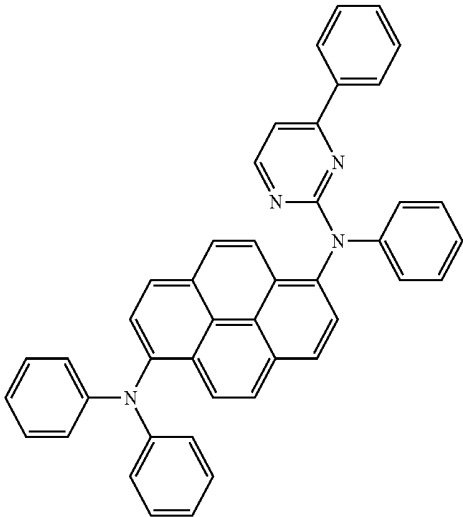


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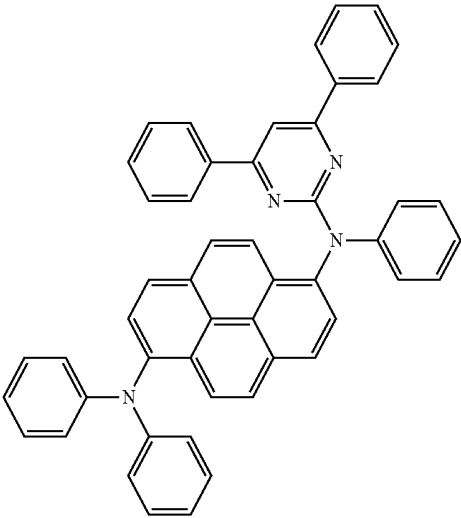
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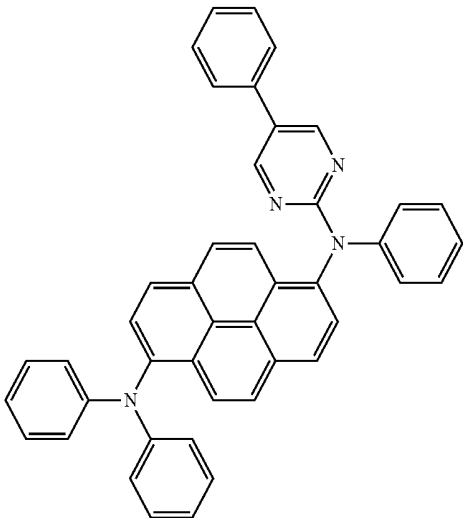
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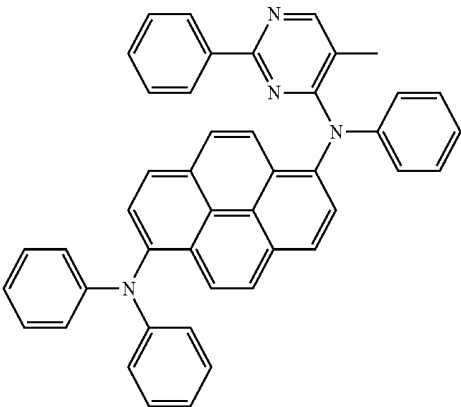
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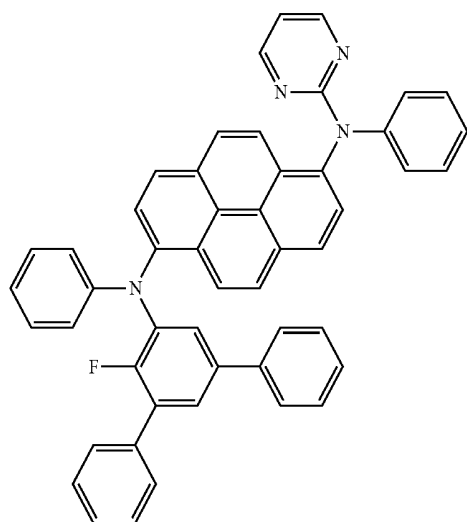


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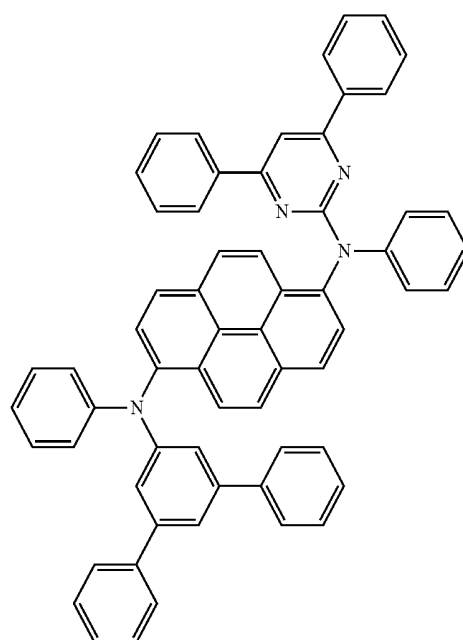
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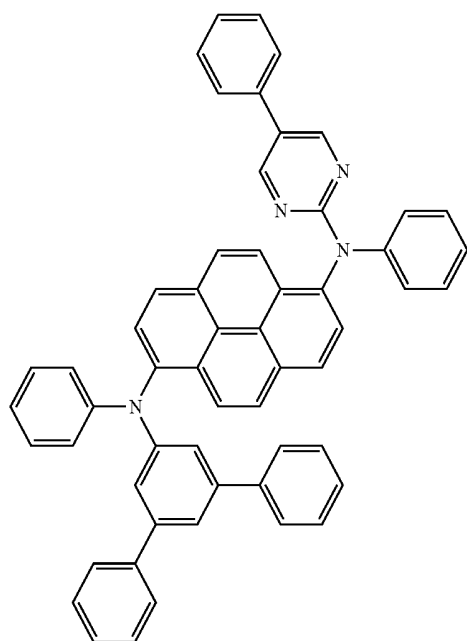
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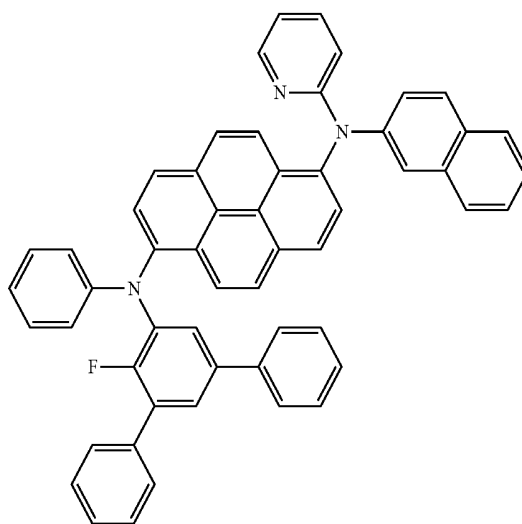


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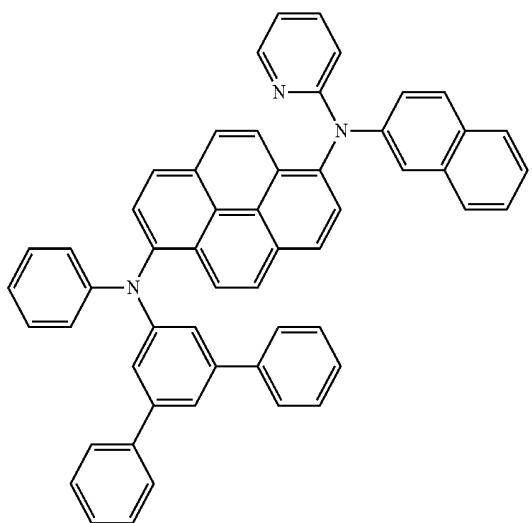


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17. An organic light-emitting diode, comprising:
a first electrode;
a second electrode facing the first electrode; and
an organic layer between the first electrode and the second electrode and including an emission layer, wherein the organic layer includes the pyrene-based compound of claim 1.

18. The organic light-emitting diode of claim 17, wherein the organic layer further includes:

a hole transportation region between the first electrode and the emission layer, the hole transportation region including at least one of a hole injection layer, a hole transportation layer, a functional layer having a hole injection capability and a hole transport capability, a buffer layer, and an electron blocking layer; and

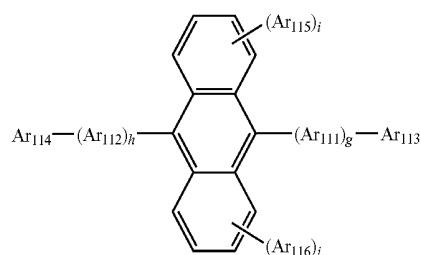
an electron transportation region between the emission layer and the second electrode, the electron transportation region including at least one of a hole blocking layer, an electron transportation layer, and an electron injection layer.

19. The organic light-emitting diode of claim 17, wherein the pyrene-based compound is:

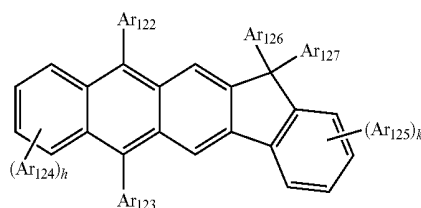
in the emission layer, and
a fluorescent dopant that emits light according to a fluorescence emission mechanism, the emission layer further including a host.

20. The organic light-emitting diode of claim 19, wherein the host includes at least one of an anthracene-based compound represented by Formula 400, below, or an anthracene-based compound represented by Formula 401, below:

<Formula 400>



<Formula 401>



wherein, in Formulae 400 and 401,

Ar₁₁₁ and Ar₁₁₂ are each independently a substituted or unsubstituted C₆-C₆₀ arylene group;

Ar₁₁₃ to Ar₁₁₆ and Ar₁₂₂ to Ar₁₂₅ are each independently a substituted or unsubstituted C₁-C₁₀ alkyl group or a substituted or unsubstituted C₆-C₆₀ aryl group;

Ar₁₂₆ and Ar₁₂₇ are each independently a C₁-C₁₀ alkyl group; and

g, h, i, j, k, and l are each independently an integer of 0 to 4.

* * * * *

专利名称(译)	基于芘的化合物和包括其的有机发光二极管		
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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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外部链接	Espacenet USPTO		

摘要(译)
基于芘的化合物和包含其的有机发光二极管，所述芘基化合物由下式1表示：

