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ABBREVIATIONS

E_{ex}	=	Excitation energy
f	=	Oscillator strength
AM1	=	Semiempirical austin model 1
B3LYP	=	Becke 's three parameter hybrid functional using the LYP Correlation functional
BH&HLYP	=	Half-and-half functionals $0.5 \cdot E_{\text{X}}^{\text{HF}} + 0.5 \cdot E_{\text{X}}^{\text{LSDA}} +$ $0.5 \cdot \Delta E_{\text{X}}^{\text{Becke88}} + E_{\text{C}}^{\text{LYP}}$
CASPT2	=	Second-order perturbation theory
CASSCF	=	Complete active space self-consistent field
CC	=	Couple cluster
CCSD	=	Singles and doubles couple cluster
CIS	=	Singles configuration interaction
CISD	=	Singles and doubles configuration interaction
CISDT	=	Singles, doubles and triples configuration interaction
CISDTQ	=	Singles, doubles, triples and quadruples configuration Interaction
(Cz) _N	=	Carbazole oligomers
(Cz-co-F) _N	=	Carbazole combined with furan oligomers
(Cz-co-Fl) _N	=	Carbazole combined with fluorine oligomers
(Cz-co-P) _N	=	Carbazole combined with phenyl oligomers
(Cz-co-Th) _N	=	Carbazole combined with thiophene oligomers
(DBZF) _N	=	Dibenzofuran oligomers
(DBZTh) _N	=	Dibenzothiophene oligomers
DFT	=	Density functional theory
E_{g}	=	Energy gap
$-\epsilon_{\text{HOMO}}$	=	Negative of HOMO energies
(FL) _N	=	Fluorene oligomers
$\Delta_{\text{H-L}}$	=	HOMO-LUMO gaps
HF	=	Hartree Fock theory
HOMO	=	Highest occupied molecular orbital
INDO/s	=	Intermediate neglect of differential overlap

LEDs	=	Light-emitting diodes
LUMO	=	Lowest unoccupied molecular orbital
MRCI	=	Multireference configuration interaction
MO	=	Molecular orbital
OLED	=	Organic light emitting diode
PA	=	Polyacetylene
PLEDs	=	Polymer light-emitting diodes
PPP	=	Poly(para-phenylene)
PPV	=	Poly(para-phenylenevinylene)
RPA	=	Random-phase approximation
RCIS	=	Restricted configuration interaction
RHF	=	Restricted Hartree-Fock
S_0	=	Ground state
S_1	=	First excited state
S_2	=	Second excited state
S_3	=	Third excited state
$S_0 \rightarrow S_1$	=	Absorption transition
$S_1 \rightarrow S_0$	=	Fluorescence transition
SAC	=	Symmetry adapted cluster
SAC-CI	=	Symmetry adapted cluster-configuration interaction
TD-DFT	=	Time-dependent density functional theory
wf	=	Wave function composition
ZINDO	=	Zerner 's intermediate neglect differential overlap