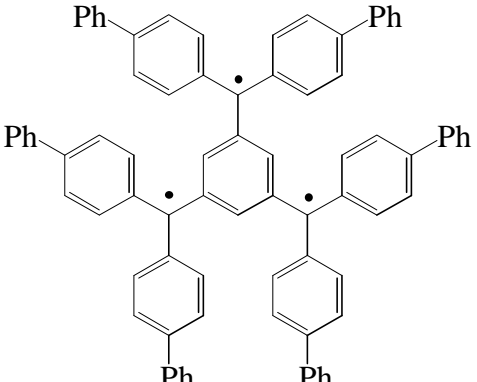
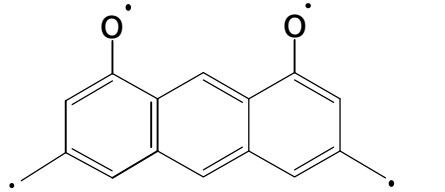
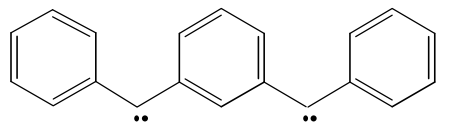
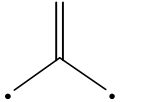
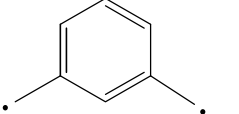
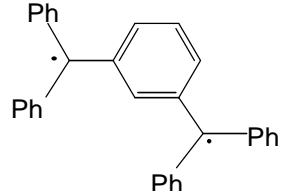
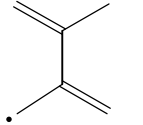
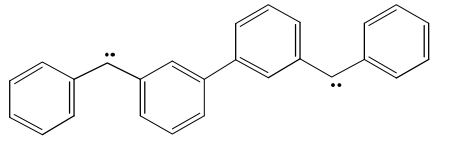
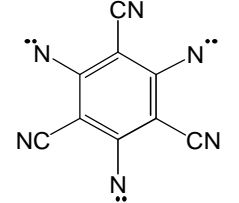
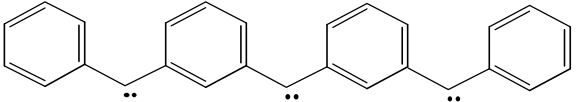
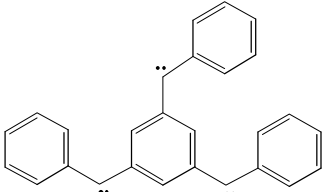
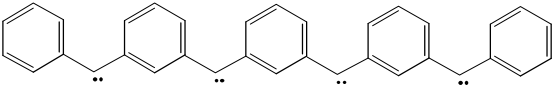
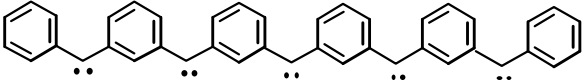
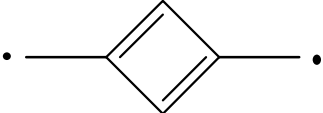
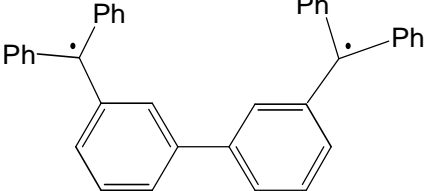
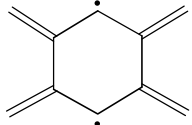
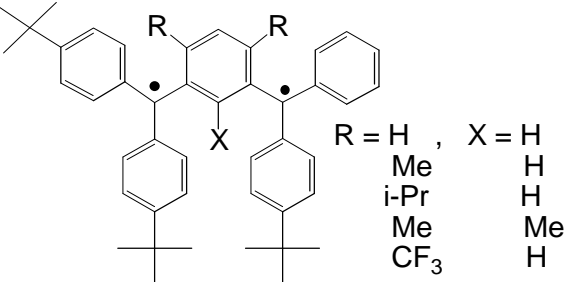
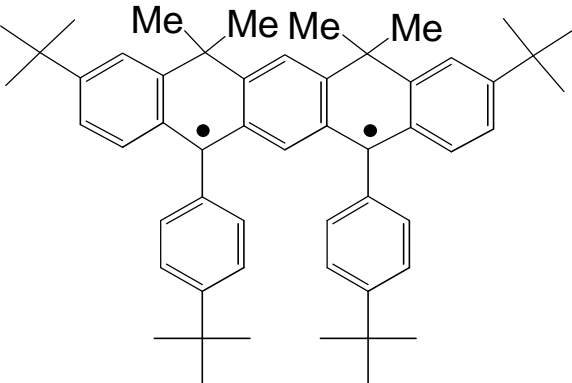
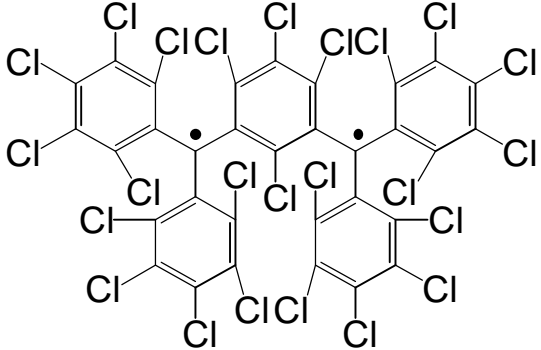
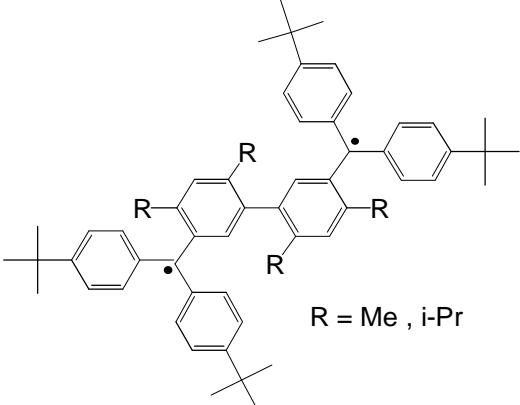
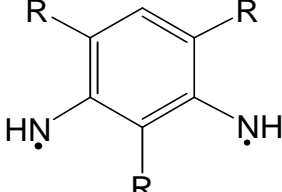
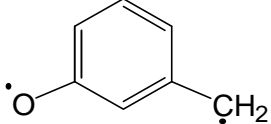
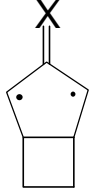
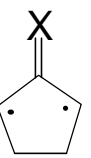
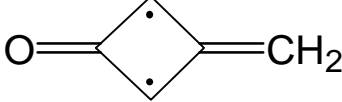
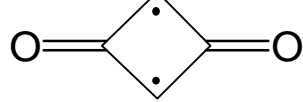
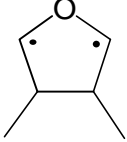
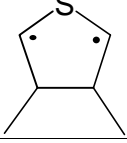
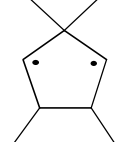
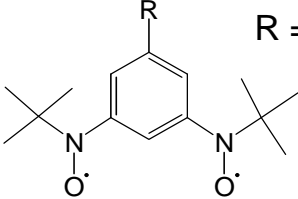
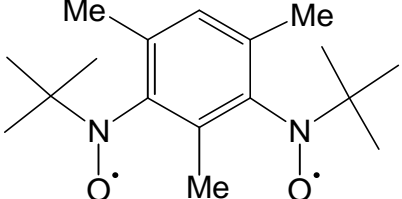
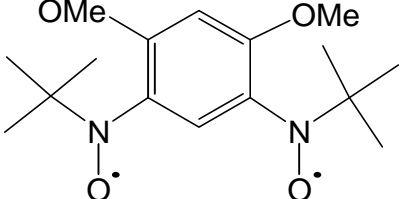
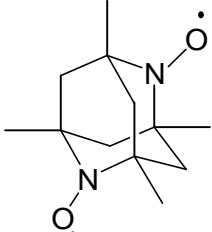
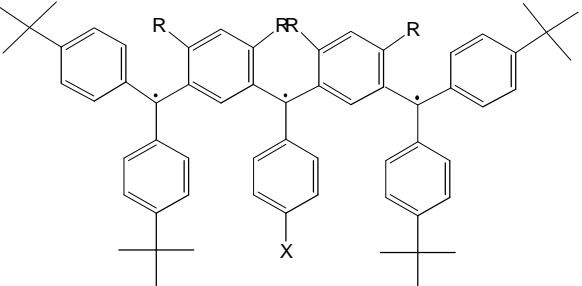
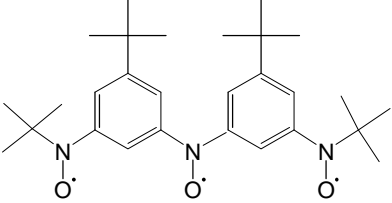
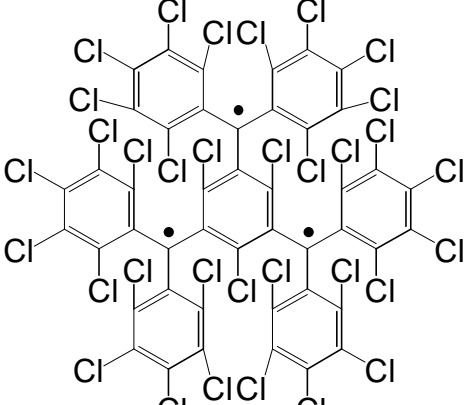
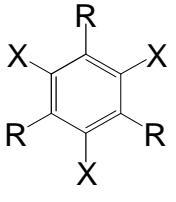
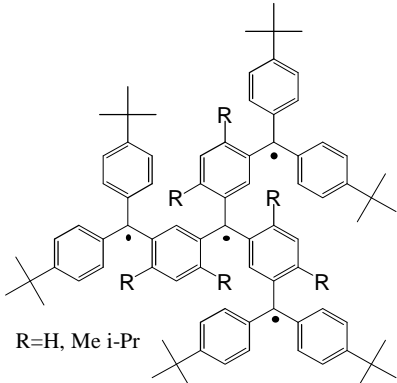
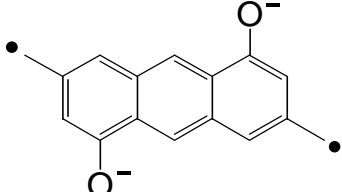
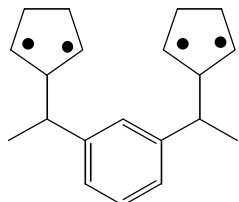
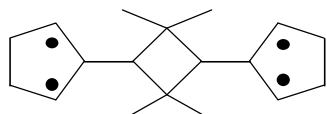


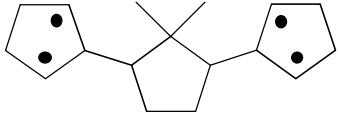
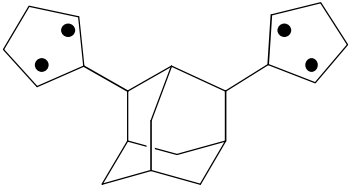
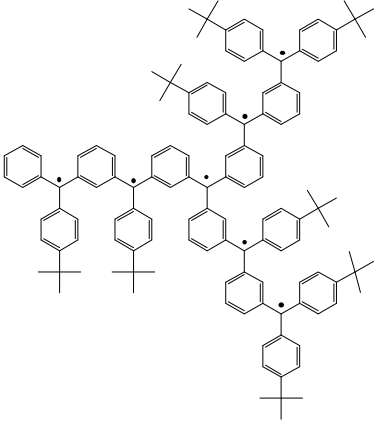
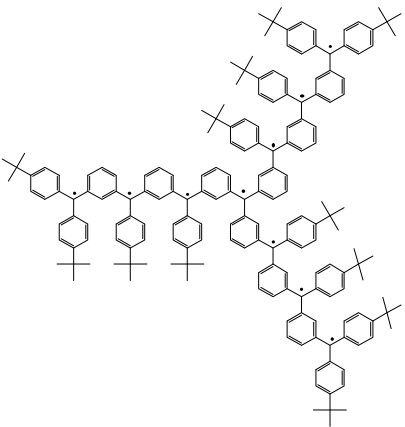
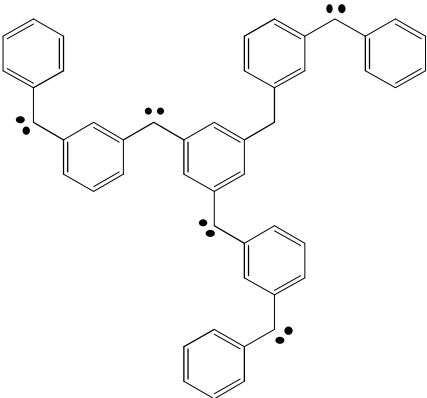
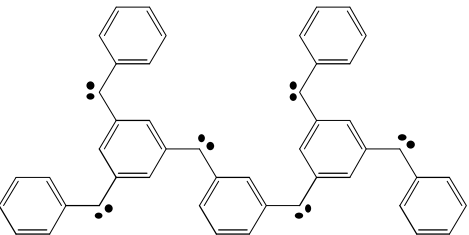
分子構造	基底状態のスピン量子数	備考	文献番号
	3/2	D=0.0041cm ⁻¹ E=0cm ⁻¹	1, 2, 3
	2	D=0.01174cm ⁻¹ E=0.00315cm ⁻¹	4, 5
	2	D=0.01731cm ⁻¹ E=0.01902cm ⁻¹	6, 7
	1		8, 9, 10
	1		11, 12, 13
	1		14, 15
	1		16~20
	0		21
	3		22

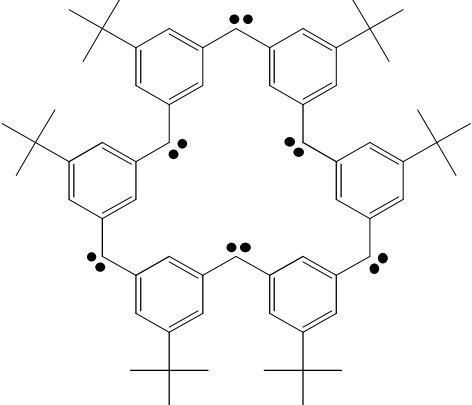
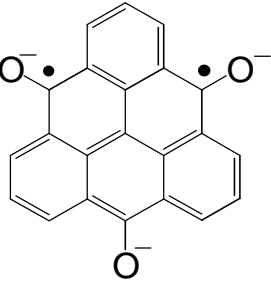
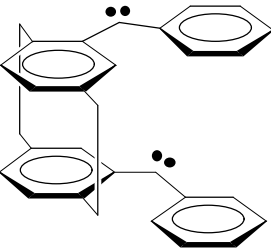
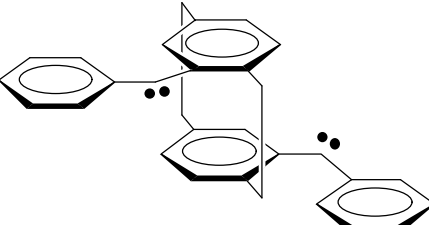
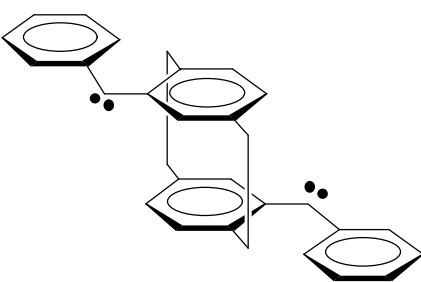
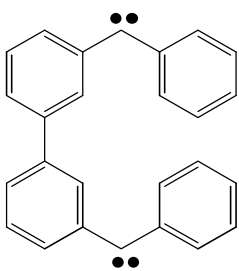
分子構造	基底状態のスピン量子数	備考	文献番号
	3	$D=0.004874\text{cm}^{-1}$ $E=0.00889\text{cm}^{-1}$	23
	3	$D=0.04158\text{cm}^{-1}$ $E=0.01626\text{cm}^{-1}$	24
	4	$D=0.03161\text{cm}^{-1}$ $E=0.00394\text{cm}^{-1}$	25
	5		26
	1		27
	1	$E_{ST}=0.3\text{kcal/mol}$	28, 29
	0		30, 31
 <p style="margin-left: 200px;"> $R = \text{H}, \text{Me}, \text{i-Pr}, \text{Me}, \text{CF}_3$ $X = \text{H}, \text{H}, \text{H}, \text{Me}, \text{H}$ </p>	1		32, 33, 34
	1		32, 33, 34

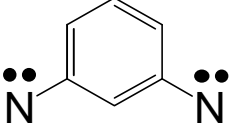
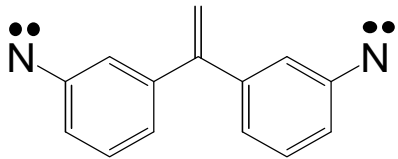
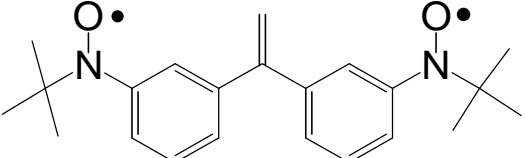
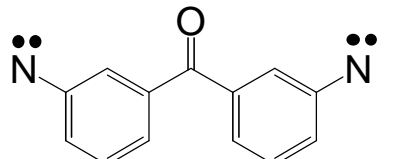
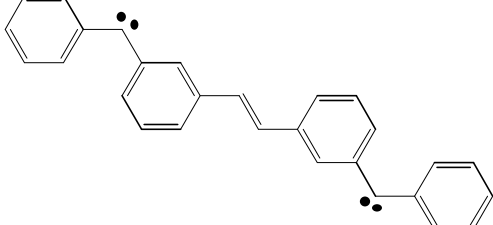
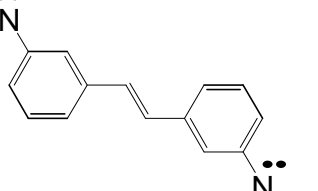
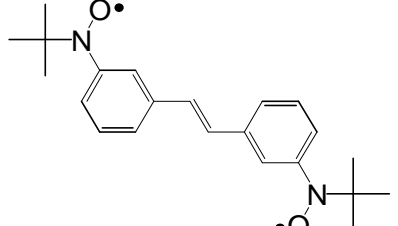
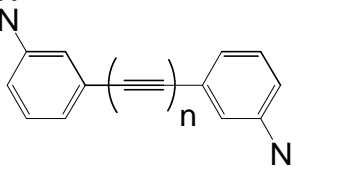
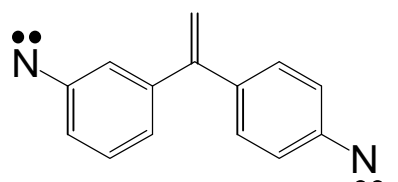
分子構造	基底状態のスピン量子数	備考	文献番号
	1		32, 33, 34
 <p>R = Me, i-Pr</p>	1	$E_{ST} =$ 0.04kcal/mol (R=Me) 0.004kcal/mol (R=i-Pr)	35
	1		36
	1		37, 38
 <p>X = C(Ph)₂, CH(Ph), C(CN)Ph, CH(CN), C(CN)₂, O</p>	0		39, 40, 41
 <p>X = C(Me)₂, C(SOMe)₂, C(SO₂Ph)₂</p>	0		39, 40, 41
	1		42
	0		43,44

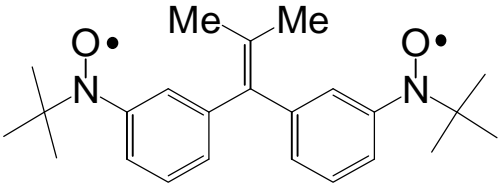
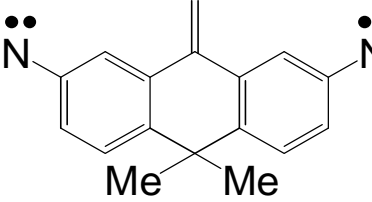
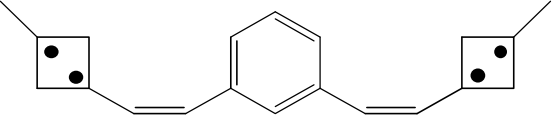
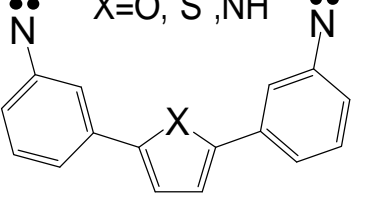
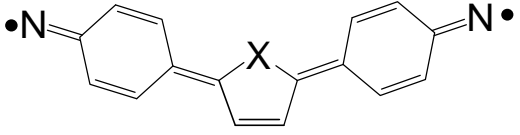
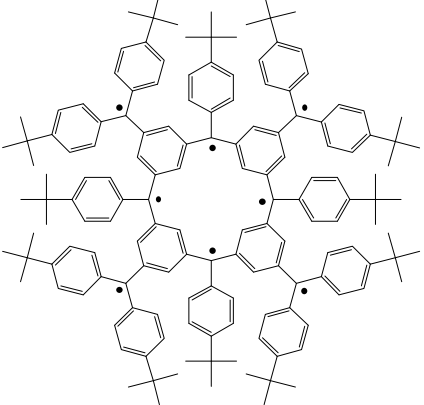
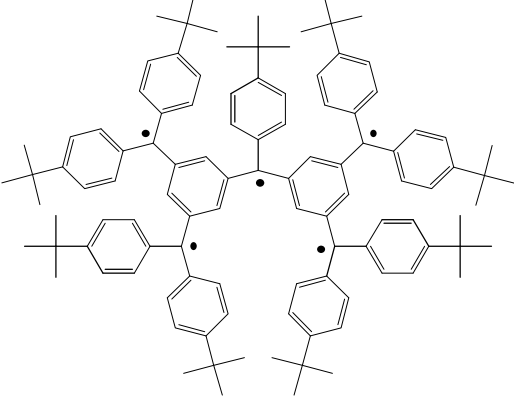
分子構造	基底状態のスピン量子数	備考	文献番号
	0		45 ~ 49
	0		45 ~ 49
	1		50, 51
 <p>R = H, t-Bu</p>	1	$E_{ST} > 1\text{kcal/mol}$	52, 53
	0		54
	0		55
	1		56, 57
 <p>R = H, Me, i-Pr, t-Bu, Ph</p> <p>X = t-Bu, t-Bu, t-Bu, Ph</p>	3/2	$E_{QD} \ 1.3\text{kcal/mol}$	58

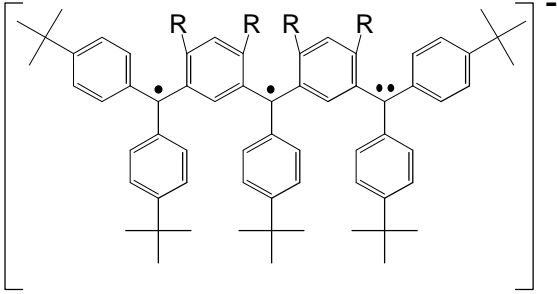
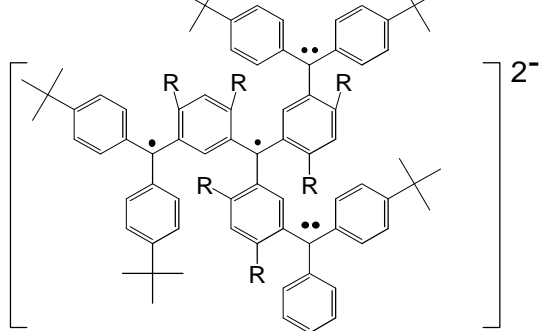
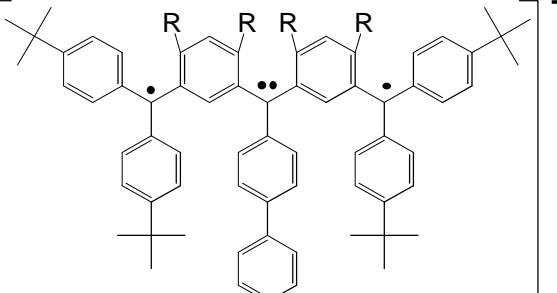
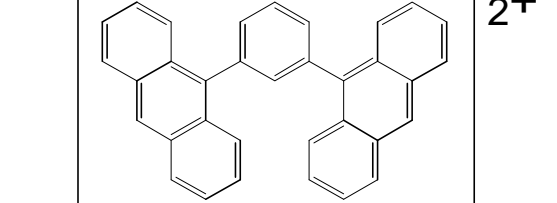
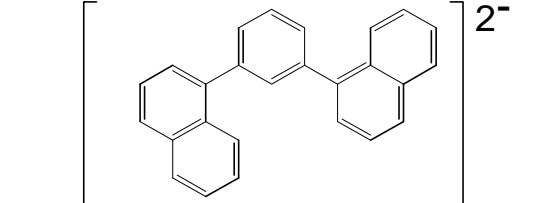
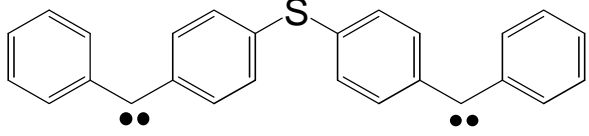
分子構造	基底状態のスピン量子数	備考	文献番号
	3/2	J 0.5kcal/mol	59
	3/2	J 0.6kcal/mol	60 ~64
 <p data-bbox="263 862 686 1086"> R=(Ph-Ph)₂C; X=H R=Ph₂N⁺; X=H R=galvinoxyl, X=H R=nitronylnitroxide, X=H R=t-butyl nitroxide, X=H, OMe R=Ph₂NN, X=CN </p>	3/2		65, 66
 <p data-bbox="167 1433 303 1467">R=H, Me i-Pr</p>	2		67 ~70
	0 or 1		71, 72
	2		73, 74, 75
	2		73, 74, 75

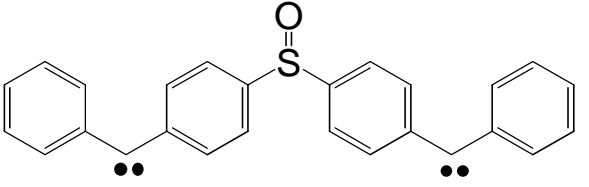
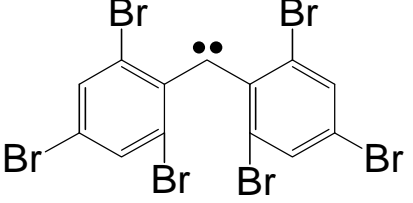
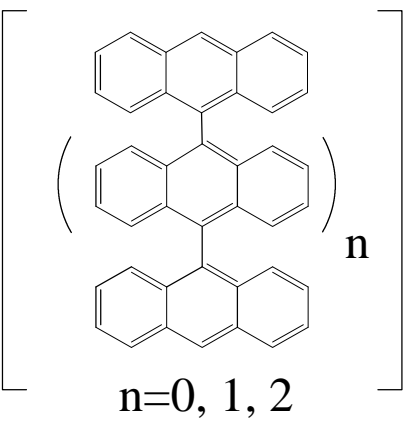
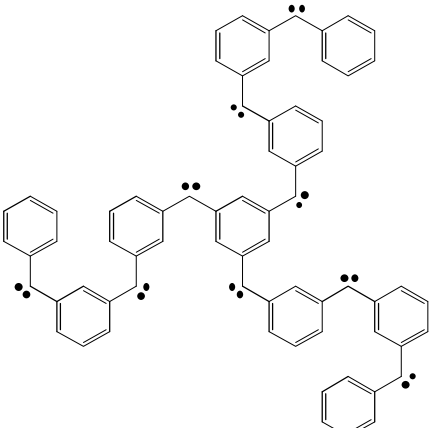
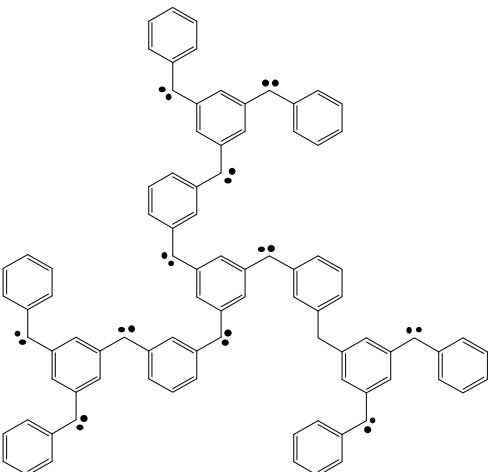
分子構造	基底状態のスピン量子数	備考	文献番号
	2		73, 74, 75
	0		73, 74, 75
	7/2		76
	5		76
	6		77,98
	6		77,98

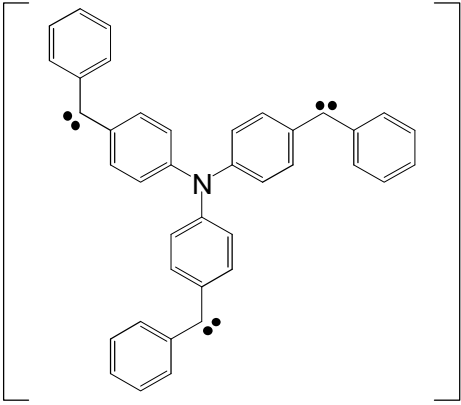
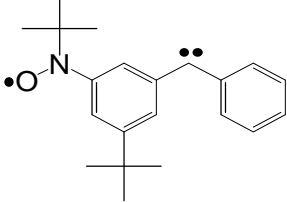
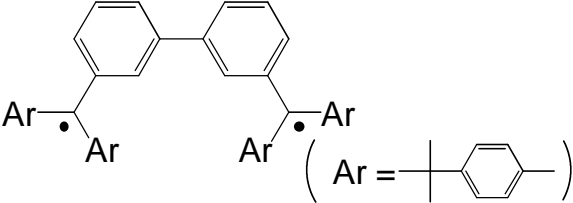
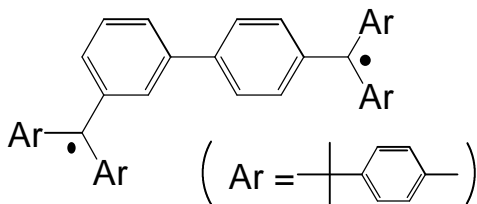
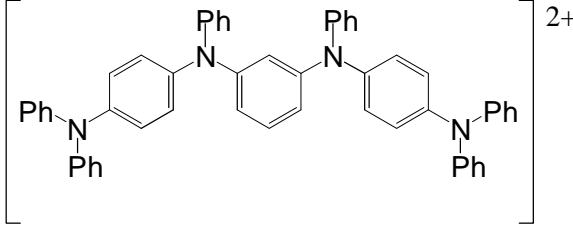
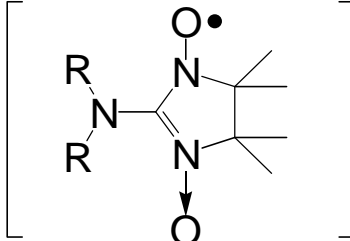
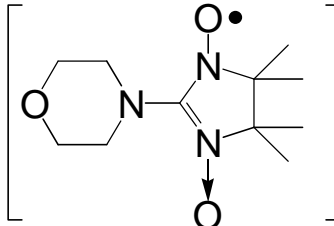
分子構造	基底状態のスピン量子数	備考	文献番号
	6		77, 98
	1	$D=0.0064\text{cm}^{-1}$ $E=0.0\text{cm}^{-1}$	78
	2	$E_{\text{QT}}=0.75\text{kJ/mol}$ (63cm^{-1})	79
	0	$E_{\text{ST}}=1.2\text{kJ/mol}$ (98cm^{-1})	79
	2	E_{ST} 60Kまでの測定	79
	0	$E=340\text{cal/mol}$ $E=770\text{cal/mol}$	80, 81

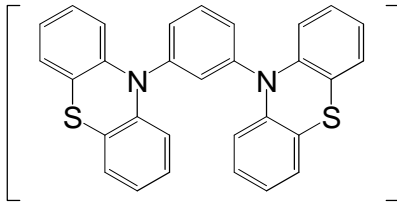
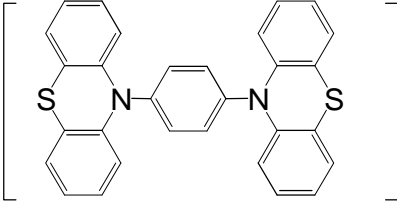
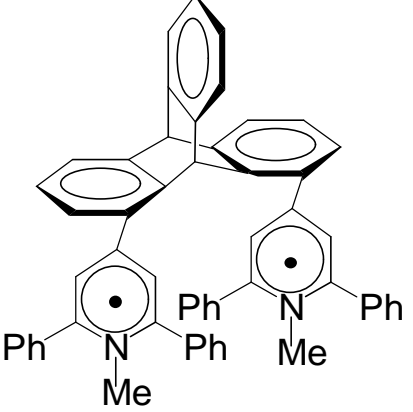
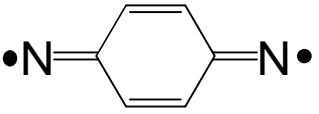
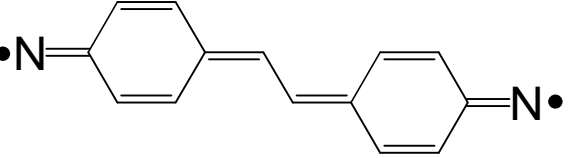
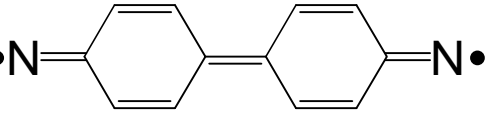
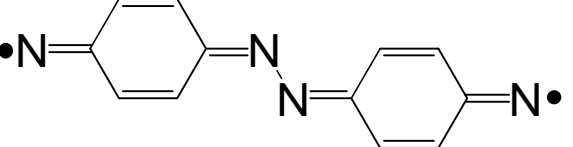
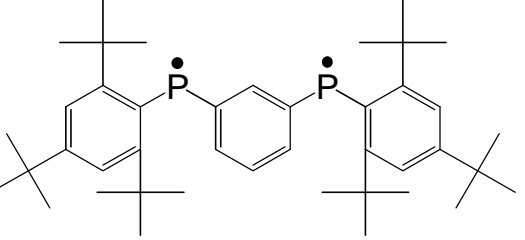
分子構造	基底状態のスピン量子数	備考	文献番号
	2		82
	0	E=75cal/mol E=85cal/mol	83, 84
	0	E=5.1cal/mol	83, 85
	0	E=48cal/mol	84
	0	E=200cal/mol	86
	0	E=240cal/mol	87
	0	E=38cal/mol	83, 85
	n=1 0 n=2 0	E=200cal/mol E=100cal/mol	88
	0	E=120cal/mol	83

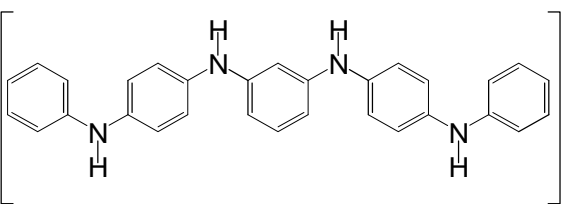
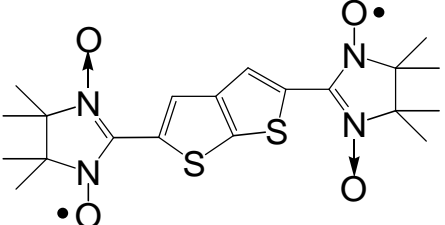
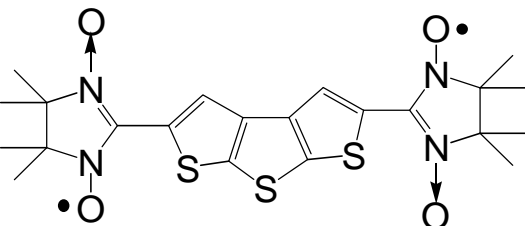
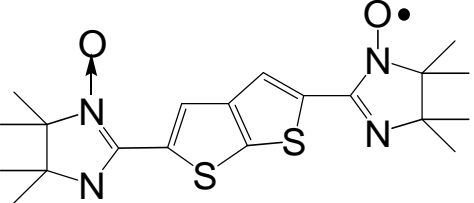
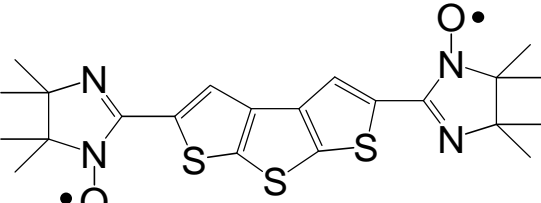
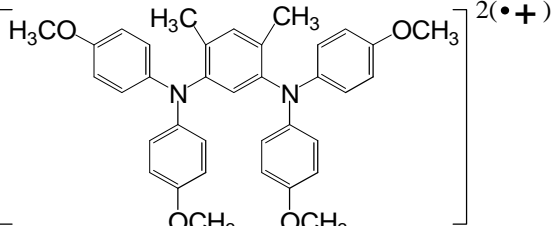
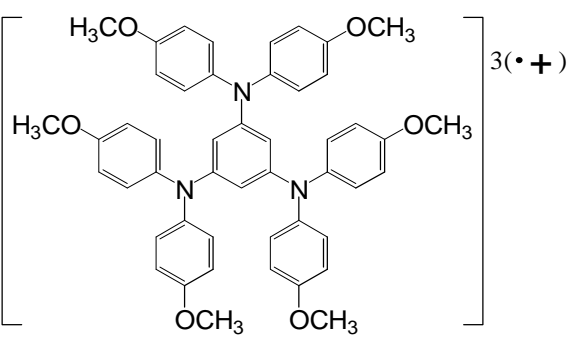
分子構造	基底状態のスピン量子数	備考	文献番号
	0	E=9.7cal/mol	83, 85
	0	$E_{ST}=810\text{J/mol}$	92
	2		91
	0		90
	0		90
	4		89
	5/2		89

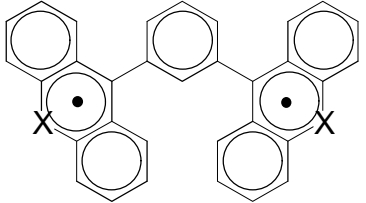
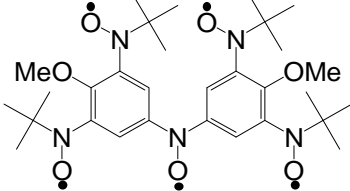
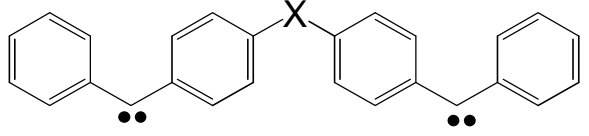
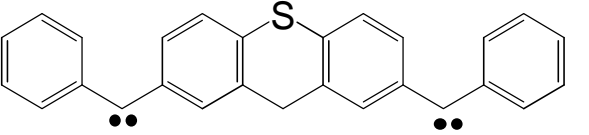
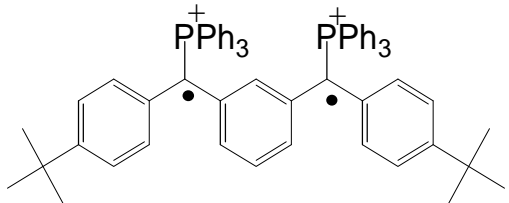
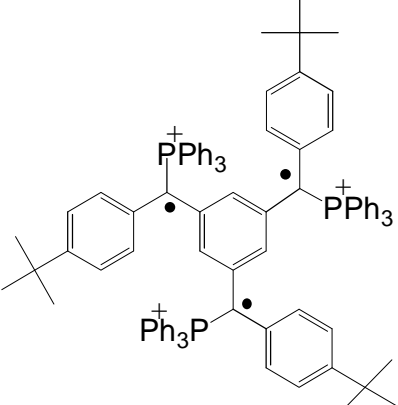
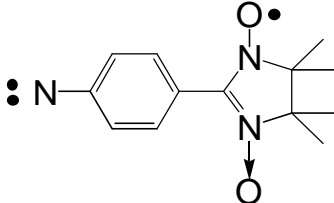
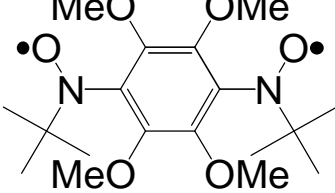
分子構造	基底状態のスピン量子数	備考	文献番号
 <p>R=H, Me, i-Pr</p>	1		94
 <p>R=H, Me, i-Pr</p>	1		94
	0		94
	1	D=0.00491cm ⁻¹	93
	1	D=0.00660cm ⁻¹	93
	2		97

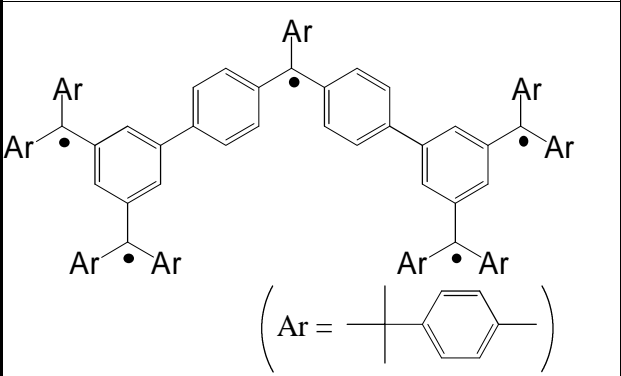
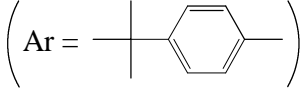
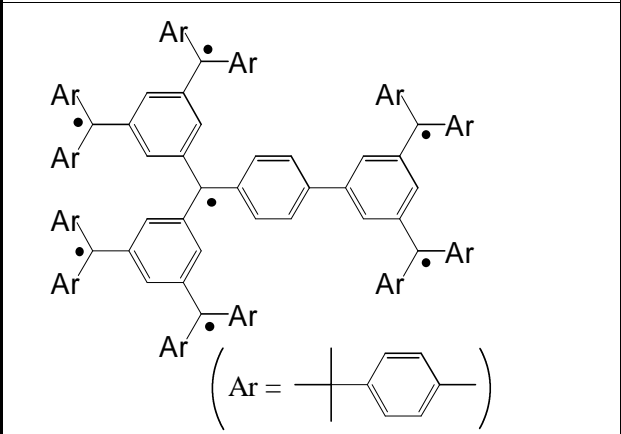
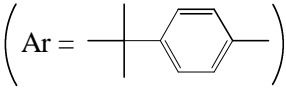
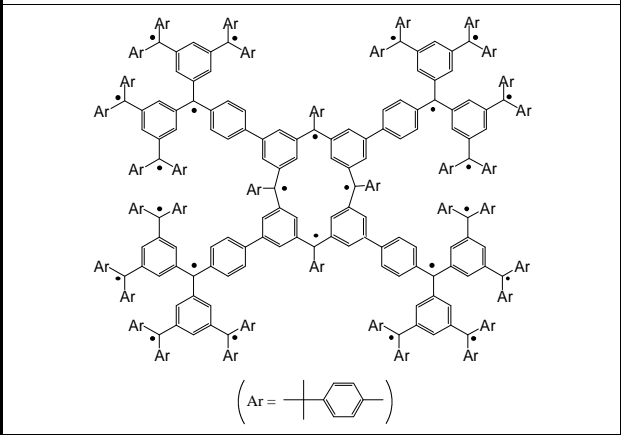
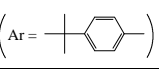
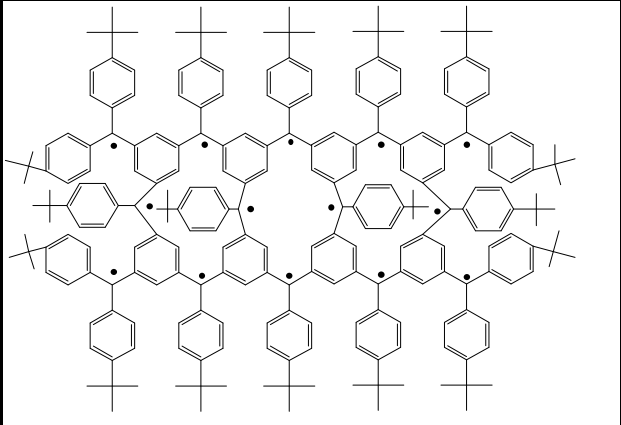
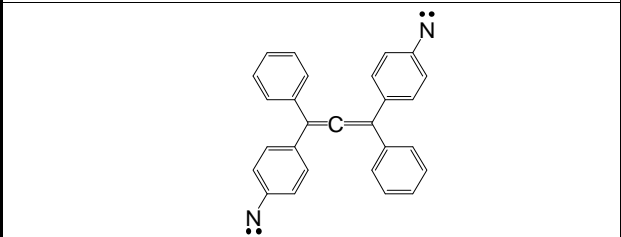
分子構造	基底状態のスピン量子数	備考	文献番号
	0		97
	1		96
	0		95
	9		102
	7 (カルベンが部分的に消滅)		100, 102

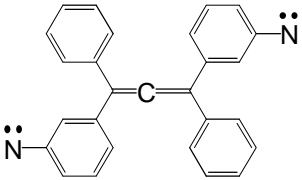
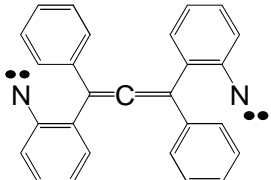
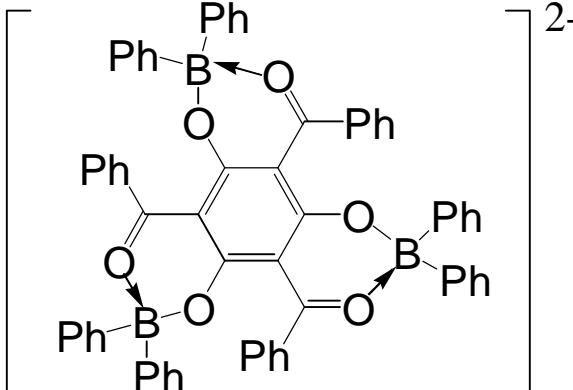
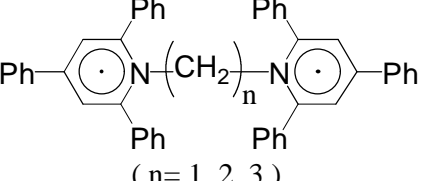
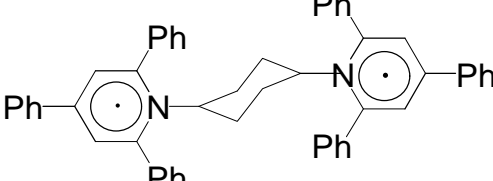
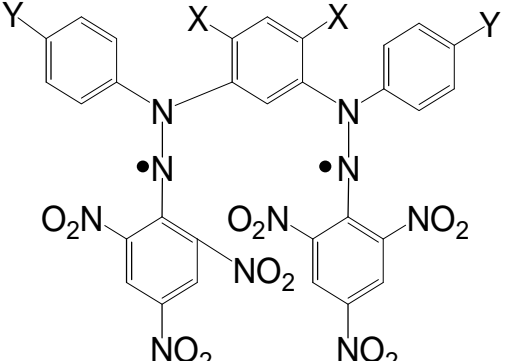
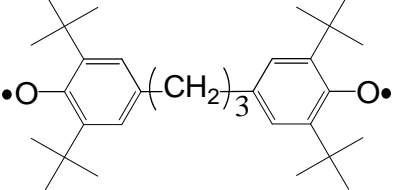
分子構造	基底状態のスピン量子数	備考	文献番号
	3/2		101
	3/2		99
	0	$E_{ST}=0.1\text{kcal/mol}$	106
	1		106
	1		105
	1		104
	1		104

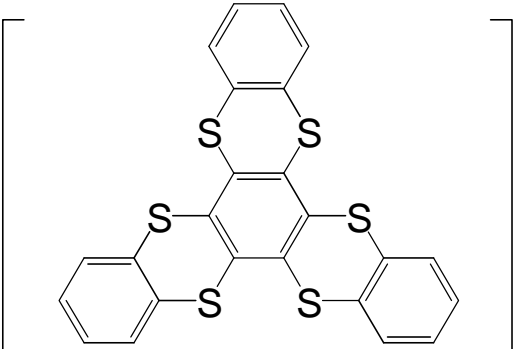
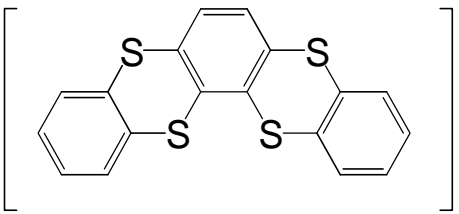
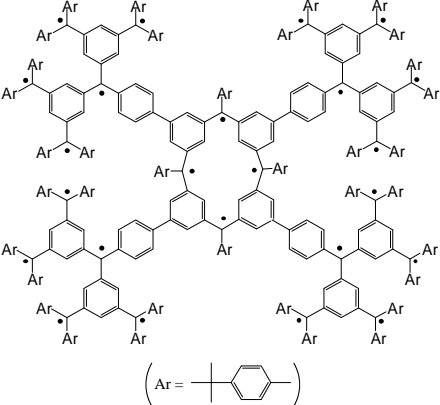
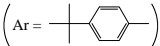
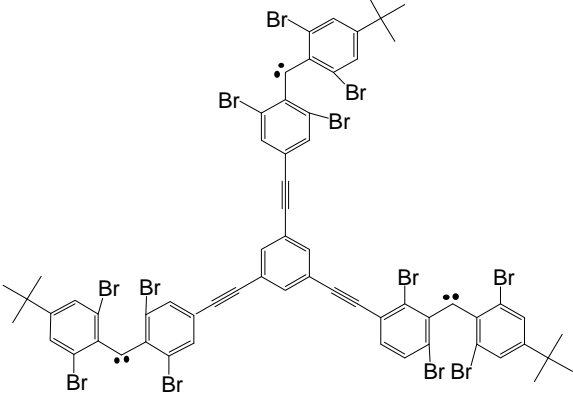
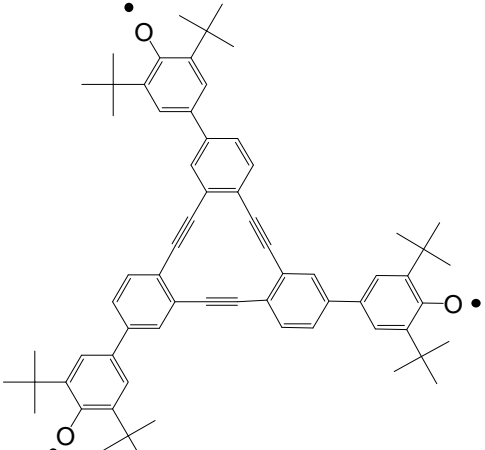
分子構造	基底状態のスピン量子数	備考	文献番号
	0	$E_{ST}=28\text{cal/mol}$	103
	1		103
	1		111
	0	$E_{ST}=0.722\text{kcal/mol}$ $D=0.169\text{cm}^{-1}$	110
	0	$E_{ST}=0.750\text{kcal/mol}$ $D=0.191\text{cm}^{-1}$	110
	0	$E_{ST}=0.472\text{kcal/mol}$ $D=0.123\text{cm}^{-1}$	110
	0	$E_{ST}=0.251\text{kcal/mol}$ $D=0.058\text{cm}^{-1}$	110
	1		109

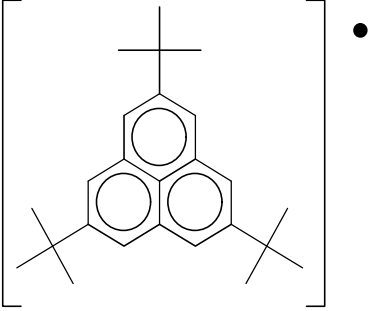
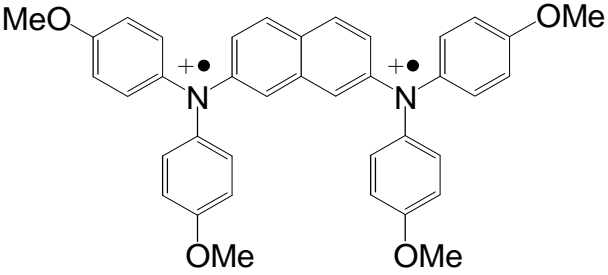
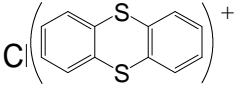
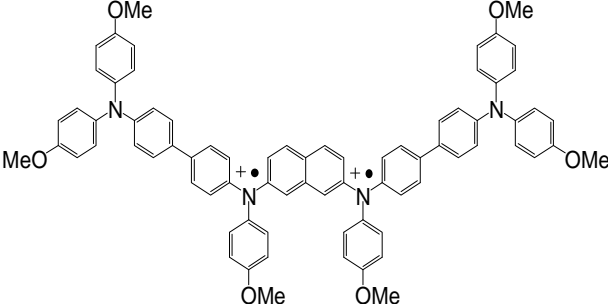
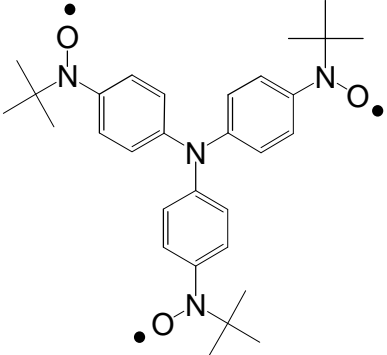
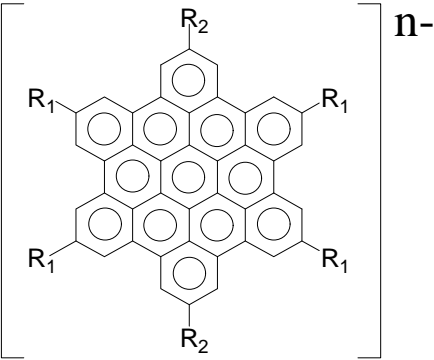
分子構造	基底状態のスピン量子数	備考	文献番号
	1	E=118MHz D=0.0MHz	108
	0		107
	0		107
	0	2J=-18K	107
	0	E _S E _T	107
	1		116
	3/2		116

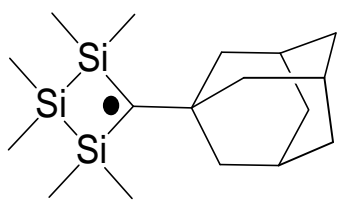
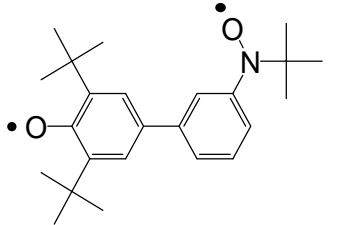
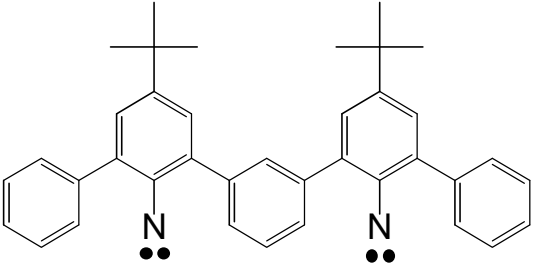
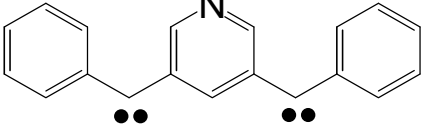
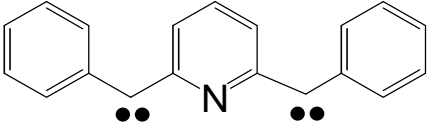
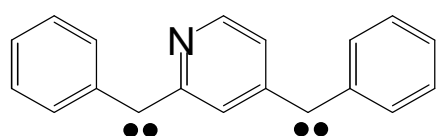
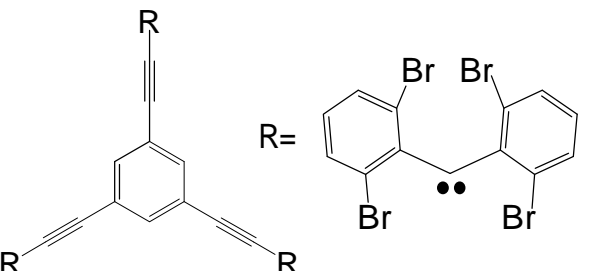
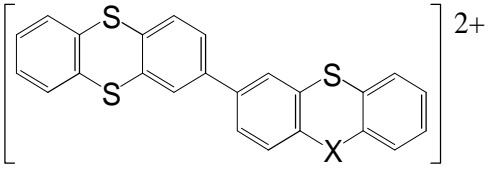
分子構造	基底状態のスピン量子数	備考	文献番号
	$X=O$ $X=S$ 1 0	$X=S$ $E_{ST}=82\text{cal/mol}$	115
	5/2		114
	$X=S$ $S=2$ $X=SO$ $S=0$ $X=SO_2$ $S=0$	$J=11\text{K}$ $J=-30\text{K}$ $J=-92\text{K}$	113
	0	$J=-21\text{K}$	113
	1		112
	3/2		112
	3/2	$ D =0.277\text{cm}^{-1}$ $ D =0.003\text{cm}^{-1}$	120
	0	$ D =0.013\text{cm}^{-1}$ $E_{ST}=14.7 \pm 0.12\text{cal/mol}$	119

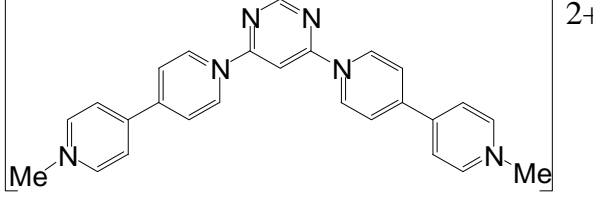
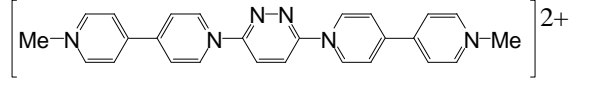
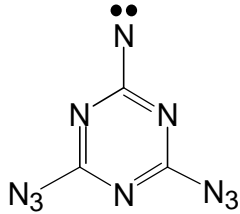
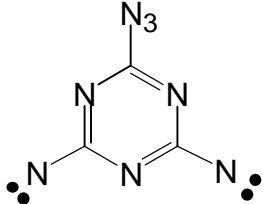
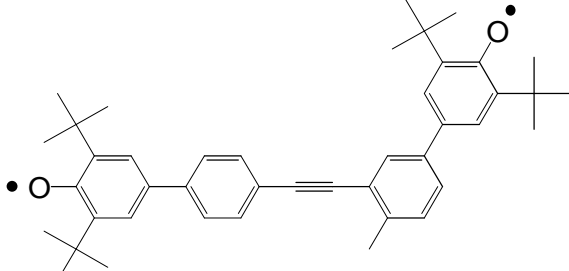
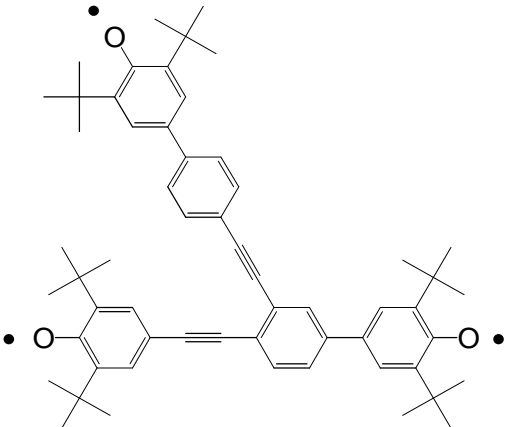
分子構造	基底状態のスピン量子数	備考	文献番号
 <p style="text-align: center;">(Ar = )</p>	2.40 (測定値)		118
 <p style="text-align: center;">(Ar = )</p>	3.28 (測定値)		118
 <p style="text-align: center;">(Ar = )</p>	7.2 (測定値)		118
	6.2 (測定値)		117
	2		124

分子構造	基底状態のスピン量子数	備考	文献番号
	2		124
	0		124
	1	$ D = 0.014 \text{ cm}^{-1}$ $ E = 0.00 \text{ cm}^{-1}$	123
 (n = 1, 2, 3)	0	$E_{\text{ST}} > 1.7 \text{ kJ/mol}$ (n=2) $E_{\text{ST}} < 230 \text{ J/mol}$ (n=1, 3)	122
	1 or (1 and 0 縮退)		122
	1 or (1 and 0 縮退)	$\left\{ \begin{array}{l} X=H, Y=H \quad (139) \\ X=H, Y=Ome \quad (139) \\ X=Me, Y=OMe(139) \\ X=Me, Y=Ph \quad (139) \\ X=Me, Y=t-Bu(121,139) \end{array} \right.$	121, 139
	1 or (1 and 0 縮退)		128

分子構造	基底状態のスピン量子数	備考	文献番号
	1	$ D = 0.012 \text{ cm}^{-1}$ $ E = 0.00 \text{ cm}^{-1}$ (127) ESR(140)	127, 140
	0	$ D = 0.012 \text{ cm}^{-1}$ $ E = 0.00 \text{ cm}^{-1}$ $E_{\text{ST}} = 35 \pm 2 \text{ J/mol}$ (127) ESR(140)	127, 140
 <p>(Ar = )</p>	10	-0.03K	126
	3		125
	3/2	磁化曲線 $S = 3/2$ (1.8 ~ 20K) $J = 34.9 \text{ K}$ (129) $2J = 48 \text{ cm}^{-1}$ $= -0.052 \text{ K}$ (143)	129, 143

分子構造	基底状態のスピン量子数	備考	文献番号
	1/2	結晶構造解析 ESR 磁気測定 $2J = -2000\text{K}$ UV-vis CV	130
	1	ESR Est=0.48eV  による酸化	131
	0 or 1	ESR Est 0	131
	1/2	結晶構造解析 ESR 磁気測定	132
 <p>1: $R_1 = R_2 = C_{12}H_{25}$ 2: $R_1 = C_{12}H_{25}$, $R_2 = t\text{-Bu}$ 3: $R_1 = R_2 = t\text{-Bu}$</p>		1^{2-} : 基底 1 重項 Est=15cal/mol 1^{3-} : 基底 2 重項 1^{4-} : 基底 1 重項 Est=20cal/mol 1^{5-} : S=1/2	133

分子構造	基底状態のスピン量子数	備考	文献番号
	1/2	結晶構造解析 ESR ENDOR	134
		ラジカル同士相互作用なし ESR UV-vis	135
	2	ESR $\begin{cases} D = 0.323 \text{ cm}^{-1} \\ E = 0.016 \text{ cm}^{-1} \end{cases}$ and $\begin{cases} D = 0.340 \text{ cm}^{-1} \\ E = 0.011 \text{ cm}^{-1} \end{cases}$	136
	2	シス体とトランス体の混合物	137
	0		137
	0		137
	3	2D-ESTN D = -0.040 cm ⁻¹ E = 0.007 cm ⁻¹	138
 <p>3²⁺: X=S 4²⁺: X=N-Me</p>	3 ²⁺ : 0 4 ²⁺ : 0 or 1		140

分子構造	基底状態のスピン量子数	備考	文献番号
	1	電解還元	141
	0	電解還元	142
	1	$ D = 1.402 \text{ cm}^{-1}$	143
	2	$ D = 0.280 \text{ cm}^{-1}$	143
	1	$2J = 31 \text{ cm}^{-1}$ $= -0.12 \text{ K}$	143
	3/2	$2J = 31 \text{ cm}^{-1}$ $= -0.16 \text{ K}$	143