

ORGANIC SUPERCONDUCTORS

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Although the vast majority of organic compounds are insulators, a small number of organic solids show considerable electrical conductivity. Some of these materials appear to be superconductors. The superconducting organics fall primarily into two groups: those containing fulvalenes (pentagonal rings containing sulfur or selenium) and those based on fullerenes, involving the nearly spherical cluster C_{60} .

The transition temperatures T_c of the fulvalene derivatives are shown in Table 1. The abbreviations of the various molecular groups are listed in Table 2 and their chemical structures are depicted in Figure 1. Most of the T_c 's are between 1 and 12 K. Several of the compounds only show superconductivity under pressure.

The fullerenes are A_3C_{60} compounds, where A represents a single or a combination of alkali atoms. The C_{60} cluster is shown in

Figure 2a, while Figure 2b illustrates how the alkali atoms fit into the A_3C_{60} molecule to form the A15 crystallographic structure. Their superconducting transition temperatures range from 8 to 31.3 K (see Table 3).

References

- Ishigura, T. and Yamaji, K., *Organic Superconductors*, Springer-Verlag, Berlin, 1990.
- Williams, Jack M. et al., *Organic Superconductors (Including Fullerenes)*, Prentice Hall, Englewood Cliffs, N.J., 1992.
- The Fullerenes*, Ed.: Krato, H. W., Fisher, J. E., and Cox, D. E., Pergamon Press, Oxford, 1993.
- Schluter, M. et al., in *The Fullerenes* (Ref. 3), p. 303.

TABLE 1. Critical Pressure and Maximum Critical Temperature of Organic Superconductors

Material	P_c /kbar	T_c /K	Material	P_c /kbar	T_c /K
(TMTSF) ₂ PF ₆	6.5	1.2	β -(ET) ₂ IBr ₂	0	2.8
(TMTSF) ₂ AsF ₆	9	1.3	β -(ET) ₂ AuI ₂	0	4.8
(TMTSF) ₂ SbF ₆	11	0.4	(ET) ₄ Hg _{2.89} Cl ₈	0	4.2
(TMTSF) ₂ TaF ₆	12	1.4	(ET) ₄ Hg _{2.89} Br ₈	12	1.8
(TMTSF) ₂ ClO ₄	0	1.4	(ET) ₃ Cl ₂ (H ₂ O) ₂	16	2
(TMTSF) ₂ ReO ₄	9.5	1.3	κ -(ET) ₂ Cu(NCS) ₂	0	10.4
(TMTSF) ₂ FSO ₃	5	3	κ -(d-ET) ₂ Cu(NCS) ₂	0	11.4
(ET) ₄ (ReO ₄) ₂	4.5	2	(DMET) ₂ Au(CN) ₂	1.5	0.9
β_L -(ET) ₂ I ₃	0	1.4	(DMET) ₂ AuI ₂	5	0.6
β_H -(ET) ₂ I ₃	0	8.1	(DMET) ₂ AuBr ₂	0	1.9
γ -(ET) ₃ I _{2.5}	0	2.5	(DMET) ₂ AuCl ₂	0	0.9
ϵ -(ET) ₂ I ₃ (I ₈) _{0.5}	0	2.5	(DMET) ₂ I ₃	0	0.6
α -(ET) ₂ I ₃ I ₂ -doped	0	3.3	(DMET) ₂ IBr ₂	0	0.7
α_t -(ET) ₂ I ₃	0	8	(MDT-TTF) ₂ AuI ₂	0	3.5
$\epsilon \rightarrow \beta$ -(ET) ₂ I ₃ ^a	0	6	TTF[Ni(dmit) ₂] ₂	2	1.6 ^b
θ -(ET) ₂ I ₃	0	3.6	TTF[Pd(dmit) ₂] ₂	20	6.5
κ -(ET) ₂ I ₃	0	3.6	(CH ₃) ₄ N[Ni(dmit) ₂] ₂	7	5

^a Converted from ϵ -type to β -type by thermal treatment.

^b For 7 kbar.

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TABLE 2. List of Symbols and Abbreviations

TTF	tetrathiafulvalene
TMTSF	tetramethyltetraselenafulvalene
BEDT-TTF or "ET"	bis(ethylenedithio)tetrathiafulvalene
MDT-TTF	methyleneedithiotetrathiafulvalene
DMET	[dimethyl(ethylenedithio)diselenadithiafulvalene]
dmit	4,5-dimercapto-1,3-dithiole-2-thione
T_c	transition temperature to superconducting state
P_c	minimum pressure required for superconducting transition

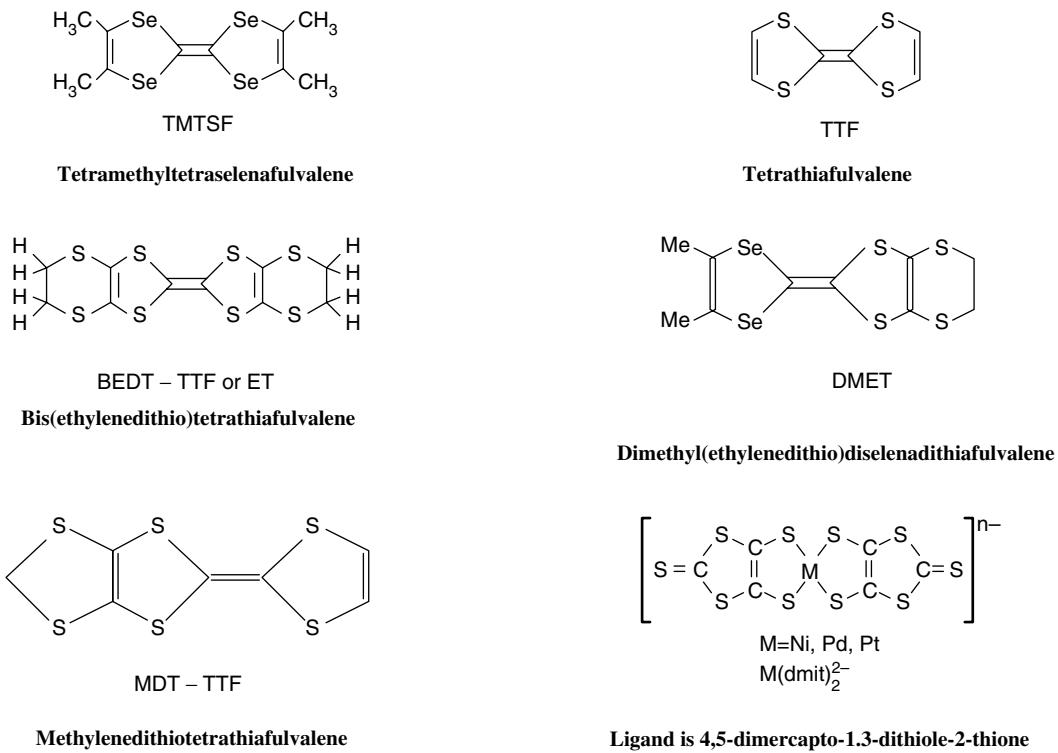


FIGURE 1. Structures of various donor molecules and acceptor species.

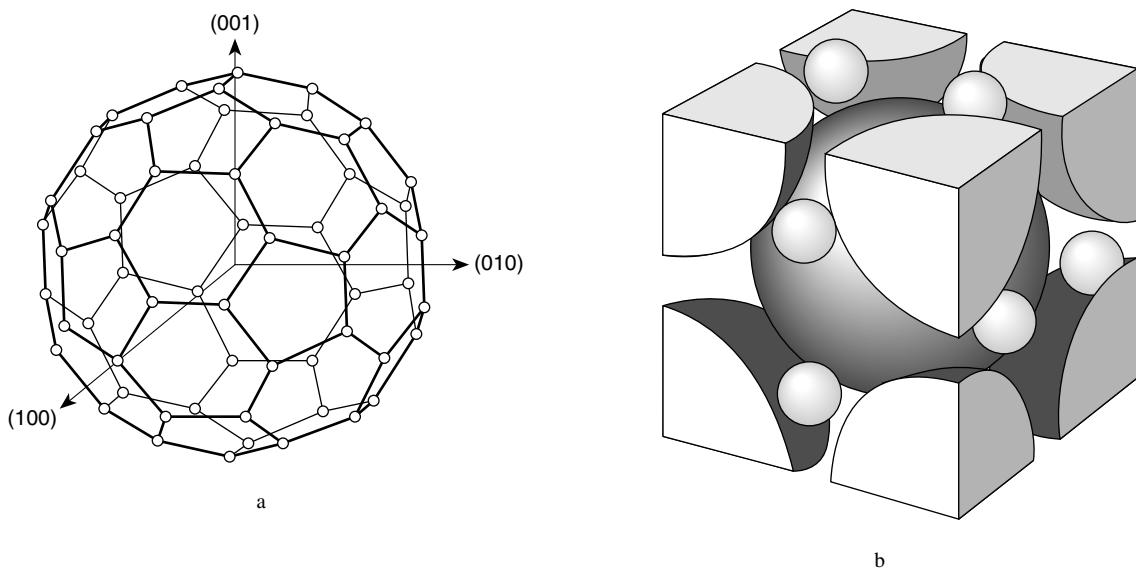


FIGURE 2. (a) C_{60} cluster placed in a fcc lattice. Each crystal axis crosses a double bond shared by two hexagons. (b) A hypothetical A_3C_{60} with the A15 structure. The structure can be seen to be an ordered defect structure of A_3C_{60} .

TABLE 3. Unit Cell and T_c for FCC-A₃C₆₀

	Lattice parameter(s) (Å)	T_c /K
Na ₂ Rb _{0.5} Cs _{0.5} C ₆₀	14.148(3)	8.0
Na ₂ CsC ₆₀ No. 1 ^a	14.132(2)	10.5
Na ₂ CsC ₆₀ No. 2 ^a	14.176(9)	14.0
K ₃ C ₆₀	14.253(3)	19.3
K ₂ RbC ₆₀	14.299(2)	21.8
Rb ₂ KC ₆₀ No. 1 ^a	14.336(1)	24.4
Rb ₂ KC ₆₀ No. 2 ^a	14.364(5)	26.4
Rb ₃ C ₆₀	14.436(2)	29.4
Rb ₂ CsC ₆₀	14.493(2)	31.3

^a Samples labeled No. 1 and No. 2 have the same nominal composition.

From Schluter, M et. al., *The Fullerenes*, Ed.: Krato, H.W., Fisher, J.E., and Cox, D.E., Pergamon Press, Oxford, 1993. With permission.