

Supplementary information.

Table 1. IR spectra of starting and referring compounds and complexes **1** and **2**.

Component	C ₇₀	(TDAE ^{•+})(C ₆₀ ^{•-}) [1]	1	2
TDAE		872w 1060m 1143w 1355m 1396s 1517m	874m 1060m 1145w 1355m 1392s* 1518m	874m 1060m 1144w 1354m 1392s* 1508m-sp 1513m-sp 1518m-sp
Fullerene	C ₇₀ 457m 534s 564m 576s 642m 673m 794m 1132w 1413w 1429s	C ₆₀ ^{•-} 514w 526w 575s 1182w 1389s	(C ₇₀ ⁻) ₂ dimer - 418w 475w 495w 508w 532s 535m 548m 557m - 568w 570w 627w 658w* 668w 671w - 712m 720m 801s 840w 902w 945w 1109w 1160w 1175m 1208w - 1244w 1257w 1279m 1307w 1391s* 1411w 1428s 1442w 1481w 1541w 1563m	(C ₇₀ ⁻) ₂ dimer 401w 417w 473w 494w 508w 530s 535m 548m 556m 560w 565w 571w 627w 658w* 667w 670w 695w 713m 720m 800s 840w 901w 944w 1109w 1158w 1175m 1209w 1229w 1244w 1257w 1279m 1307w 1392s* 1411w 1428s 1440w 1480w 1541w 1562m
C ₆ H ₄ Cl ₂			658w* 749s 1033m - 1456w	658w* 750s 1033m 1125w 1455w

sp- splitting; w-weak; m-middle; s-strong.
a- the bands of two components coincide.

[1] K. Pokhodnia, J. Papavassiliou, P. Umek, A. Omerzu, D. Mihailovič, *J. Phys. Chem.*, 1999, **110**, 3606.