Supplementary Material (ESI) for Chemical Communications

Supporting Information for

$C_6N_7Cl[N(C_6H_5)_2]_2 \cdot p\text{-}C_6H_4Me_2N$ – A mono-chloro-tri-*s*-triazine adduct

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1. Characterization

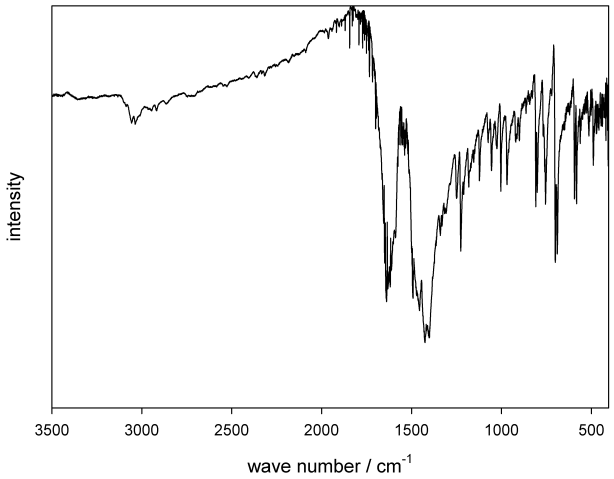

Fig. S1 FTIR spectrum of **2**·*p*-$C_6H_4Me_2$.

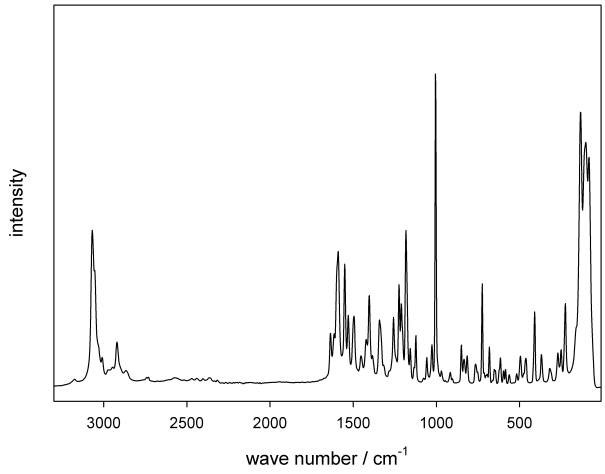

Fig. S2 Raman spectrum of **2**·*p*-$C_6H_4Me_2$.

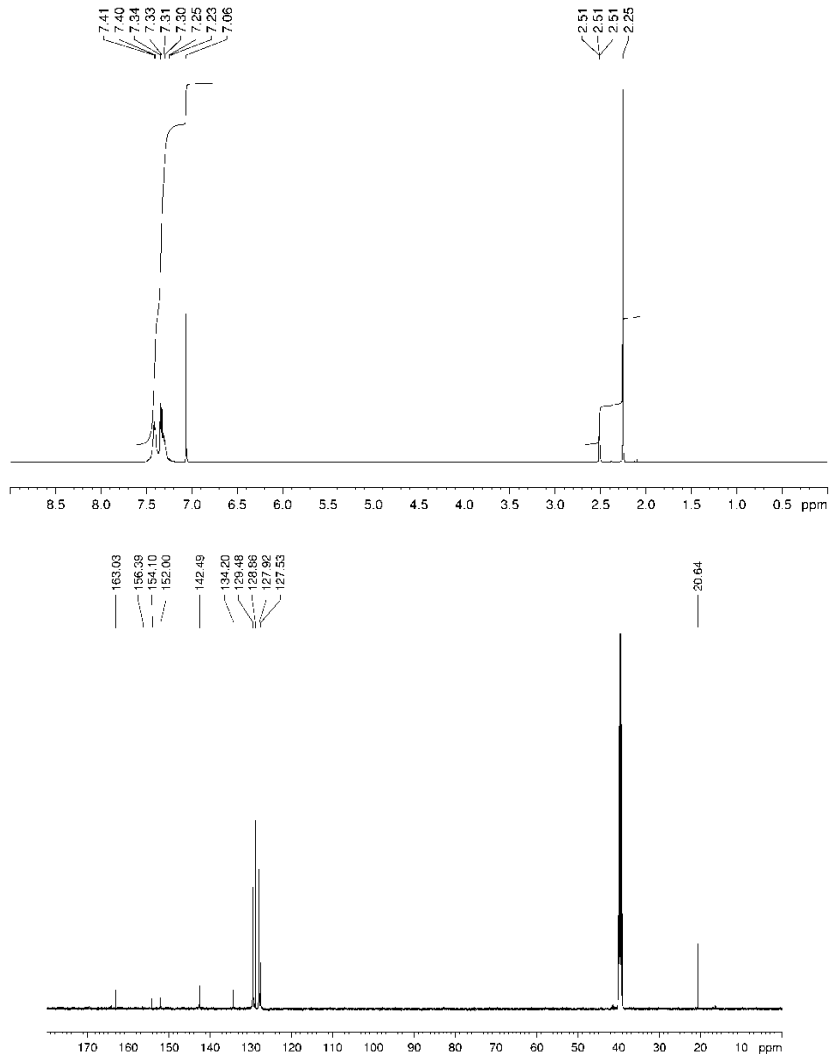

Fig. S3 ^{1}H- and ^{13}C-NMR spectra of **2·*p*-$C_6H_4Me_2$** (solvent: d_6-DMSO).

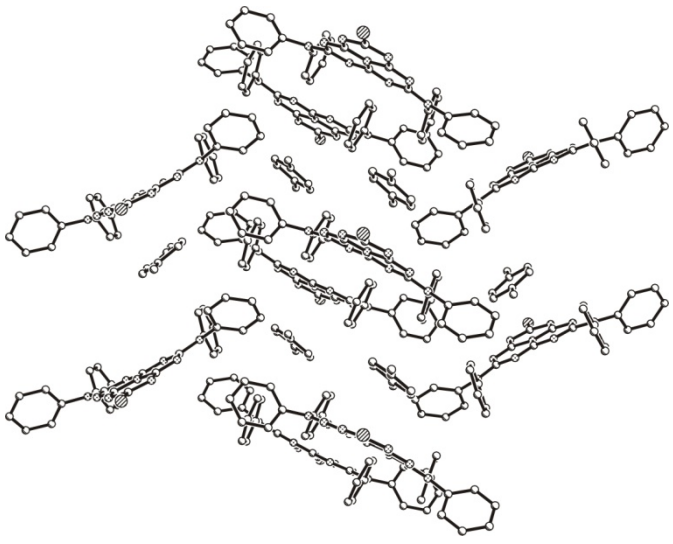

Fig. S4 Packing diagram of **2·*p*-$C_6H_4Me_2$** along the crystallographic *a* axis. Hydrogen atoms are omitted for clarity.

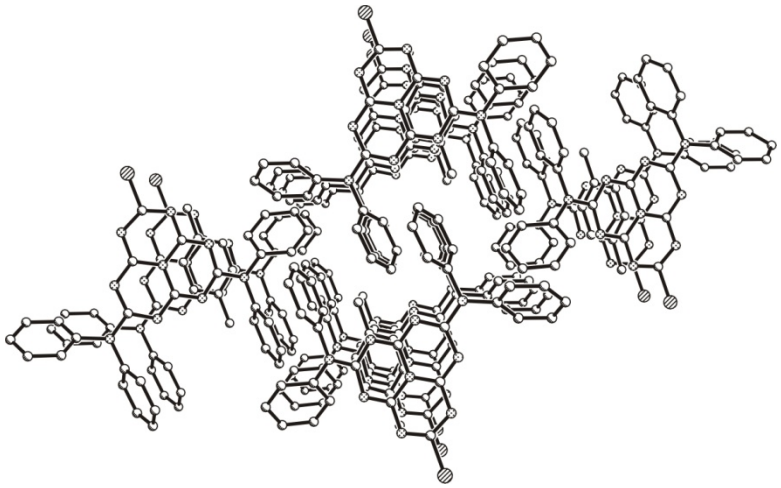

Fig. S5 Packing diagram of **2·*p*-$C_6H_4Me_2$** along the crystallographic *b* axis. Hydrogen atoms are omitted for clarity.

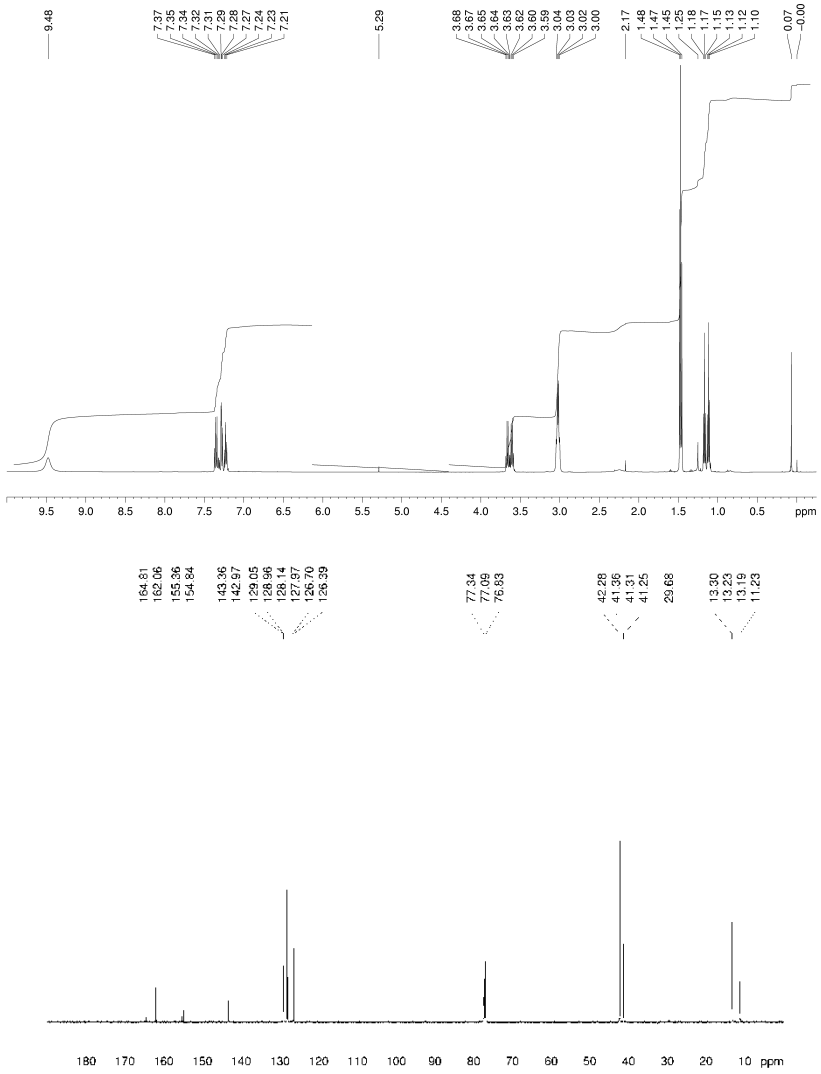

Fig. S6 ^{1}H- and ^{13}C-NMR spectra of 1,5-Bis(diphenylamino)-8-diethylamino-tri-*s*-triazin and diethylamine hydrochloride.

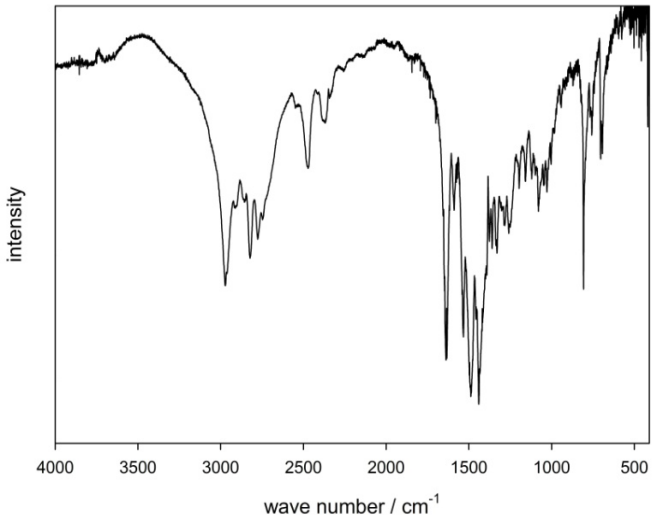

Fig. S7 FTIR spectrum of 1,5-Bis(diphenylamino)-8-diethylamino-tri-*s*-triazin.