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4-(5-Phenyl-3-trifluoromethyl-1*H*-pyrazol-1-yl)benzenesulfonamide

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Abstract: Significant twists between the aromatic rings are evident in the structure of the title compound, $C_{16}H_{12}F_3N_3O_2S$. With reference to the pyrazole plane, the N- and C-bound benzene rings form dihedral angles of 57.12 (11) and 29.75 (11)°, respectively. The dihedral angle between the benzene rings is 52.82 (11)°. The presence of N-H---O(sulfonamide) and N-H---N(pyrazole) hydrogen bonds lead to supramolecular tubes along the *b*-axis direction. These are connected into layers *via* C-H---O interactions involving a bifurcated O atom (not involved in the N-H---O hydrogen bonding). Layers stack along the *a*-axis direction.