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4-(5-Phenyl-3-trifluoromethyl-1H-pyrazol-1-yl)benzenesulfonamide
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Abstract: Significant twists between the aromatic rings are evident in the structure of the title compound, $\mathrm{C}_{16} \mathrm{H}_{12} \mathrm{~F}_{3} \mathrm{~N}_{3} \mathrm{O}_{2} \mathrm{~S}$. With reference to the pyrazole plane, the N - and C-bound benzene rings form dihedral angles of 57.12 (11) and 29.75 (11) ${ }^{\circ}$, respectively. The dihedral angle between the benzene rings is $52.82(11)^{\circ}$. The presence of $\mathrm{N}-\mathrm{H} . . \mathrm{O}$ (sulfonamide) and $\mathrm{N}-\mathrm{H}$ $\ldots \mathrm{N}$ (pyrazole) hydrogen bonds lead to supramolecular tubes along the $b$-axis direction. These are connected into layers via C $\mathrm{H}_{\ldots . .} \mathrm{O}$ interactions involving a bifurcated O atom (not involved in the $\mathrm{N}-\mathrm{H}_{\ldots} . \mathrm{O}$ hydrogen bonding). Layers stack along the $a$ axis direction.

