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***N*-(2-Bromophenyl)-4-methyl-*N*-(4-methylphenyl)sulfonylbenzenesulfonamide**

**M. N. Arshad, I. U. Khan, K. T. Holman, A. M. Asiri and H. M. Rafique**

**Abstract:** In the title compound, C<sub>20</sub>H<sub>18</sub>BrNO<sub>4</sub>S<sub>2</sub>, the mean planes formed by the toluene substituents are inclined at a dihedral angle of 45.34 (8)°. The bromobenzene group is disordered over two positions with an occupancy ratio of 0.74:0.26, resulting in two conformations of the ring; the two rings are oriented at a dihedral angle of 6.6 (6)° with each other. In the crystal structure, weak C-H...O interactions connect the molecules in a zigzag manner along the *a* axis.