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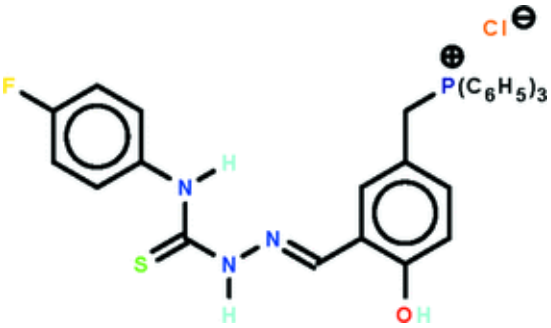

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[3-({(*E*)-2-[(4-Fluorophenyl)carbamothioyl]hydrazinylidene}methyl)-4-hydroxybenzyl]methyltriphenylphosphonium chloride

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Abstract: The cation in the title salt, $C_{33}H_{28}FN_3OPS^+\cdot Cl^-$, is highly twisted with the phosphonium group occupying a position almost normal to the central hydroxylbenzene ring [P-C-C-C tosrsion angle = -100.9 (3)°], and with the hydrazone substituent twisted out of the plane [C-C-C-N torsion angle = 13.1 (4)°]. The fluorobenzene ring is twisted out of the plane of the adjacent thiourea residue, forming a dihedral angle of 51.69 (10)°. The configuration about the C=N bond [1.281 (4) Å] is *E*, the O-H and N-H hydrogen atoms are *syn*, and in the thiourea residue, the N-H hydrogen atoms are *anti*, allowing for the formation of an intramolecular N-H...N hydrogen bond. In the crystal, dimeric aggregates mediated by N-H...S bonds are formed, which are linked to the Cl^- anions by O-H...Cl hydrogen bonds. The four-component aggregates are linked into a three-dimensional structure by C-H...Cl interactions.