IUPAC Name	(2~{S})-3-(1-methylindol-3-yl)-2-[(2-methylpropan-2-yl)oxycarbonylamino]propanoic acid
Synonyms	(2S)-2-{[(tert-butoxy)carbonyl]amino}-3-(1-methyl-1H-indol-3-yl)propanoic acid; L-Tryptophan, N-[(1,1-dimethylethoxy)carbonyl]-1-methyl
GLB Cat#	
SMILES	CC(C)(C)OC(=O)N[C@@H](CC1=CN(C2=CC=CC=C21)C)C(=O)O
Molecular Weight	318.37
Specific Rotation[a]20/D	N/A
Optical purity	Not tested
Melting Point	N/A
Solubility	Not tested
Loss on drying	≤1.0%(80°C,2h)
Water Content (K.F.)	≤2.0%
NMR Spectrum	Complies with the structure
Mass Spectrum	Complies with the structure
IR Spectrum	Complies with the structure

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