

IUPAC Name	(2R)-2-amino-3,3-dimethylbutan-1-ol
Synonyms	(R)-TERT-LEUCINOL; (2R)-2-amino-3,3-dimethylbutan-1-ol
GLB Cat#	37513
SMILES	<chem>CC(C)(C)[C@H](CO)N</chem>
Molecular Weight	117.19
Specific Rotation $[\alpha]_{20/D}$	-38.8°±1°(C=1 in EtOH)
Optical purity	N/A
Melting Point	Need testing
Solubility	Need testing
Loss on drying	Need testing
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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