3,5-dihydroxytoluene, monohydrate

Inchi: InChI=1S/C7H8O2.H2O/c1-5-2-6(8)4-7(9)3-5;/h2-4,8-9H,1H3;1H2
InchiKey: NBKPNAMTHBIMLA-UHFFFAOYSA-N
Formula: C7H10O3
SMILES: Cc1cc(O)cc(O)c1.O
Mol. weight [g/mol]: 142.15
CAS: 6153-39-5

Physical Properties

Property code	Value	Unit	Source
tf	328.00 ± 0.20	K	NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C6153395&Units=SI

Legend

tf: Normal melting (fusion) point