Molecular Origami of TeBr2

given information

ElementNames	[(Te) (Br) (Br)]	
dotted	F	
distance	251.024	Te^{1}-Br^{2}
distance	251.100	Te^{1}-Br^{1}
angle	97.969	Br^{2}-Te^{1}-Br^{1}
	378.9	Br^{2}-Br^{1}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

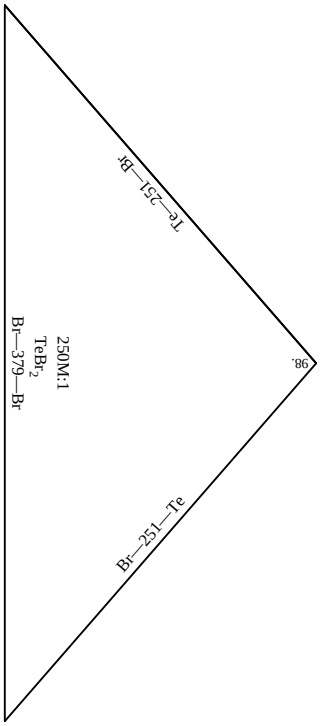
Molecular Origami of TeBr2

!Te1
Br1
Br2
TeBr2

bent

scale 250,000,000 : 1
units: pm
offsetx -1.9 offsety 0.9

View -1



Current: (centerx 2.40) (centery 5.90) (scale 250)
%%BoundingBox: 163 281 300 569 actual: 173 291 290 559 center: 231 425 actual size: 117 268
Better: (centerx 3.44) (centery 5.50) (scale 250)
%%BoundingBox: 101 317 238 605 actual: 111 327 228 595 center: 169 461 actual size: 117 268