Molecular Origami of CHBr3

given information

ElementNames	[(C) (Br) (Br) (Br) (H)]	
distance	96.988	C^1-H^1
distance	191.631	C^1-Br^3
distance	191.817	C^1-Br^1
distance	193.849	C^1-Br^2
angle	108.209	H^1-C^1-Br^1
	240.5	H^1-Br^1
angle	108.421	H^1-C^1-Br^3
	240.6	H^1-Br^3
angle	108.795	H^1-C^1-Br^2
	243.1	H^1-Br^2
angle	110.147	Br^3-C^1-Br^2
	316.1	Br^3-Br^2
angle	110.215	Br^2-C^1-Br^1
	316.3	Br^2-Br^1
angle	110.985	Br^3-C^1-Br^1
	316.	Br^3-Br^1
dopage	T	
AutoAlign	F	

structure type: XABCD

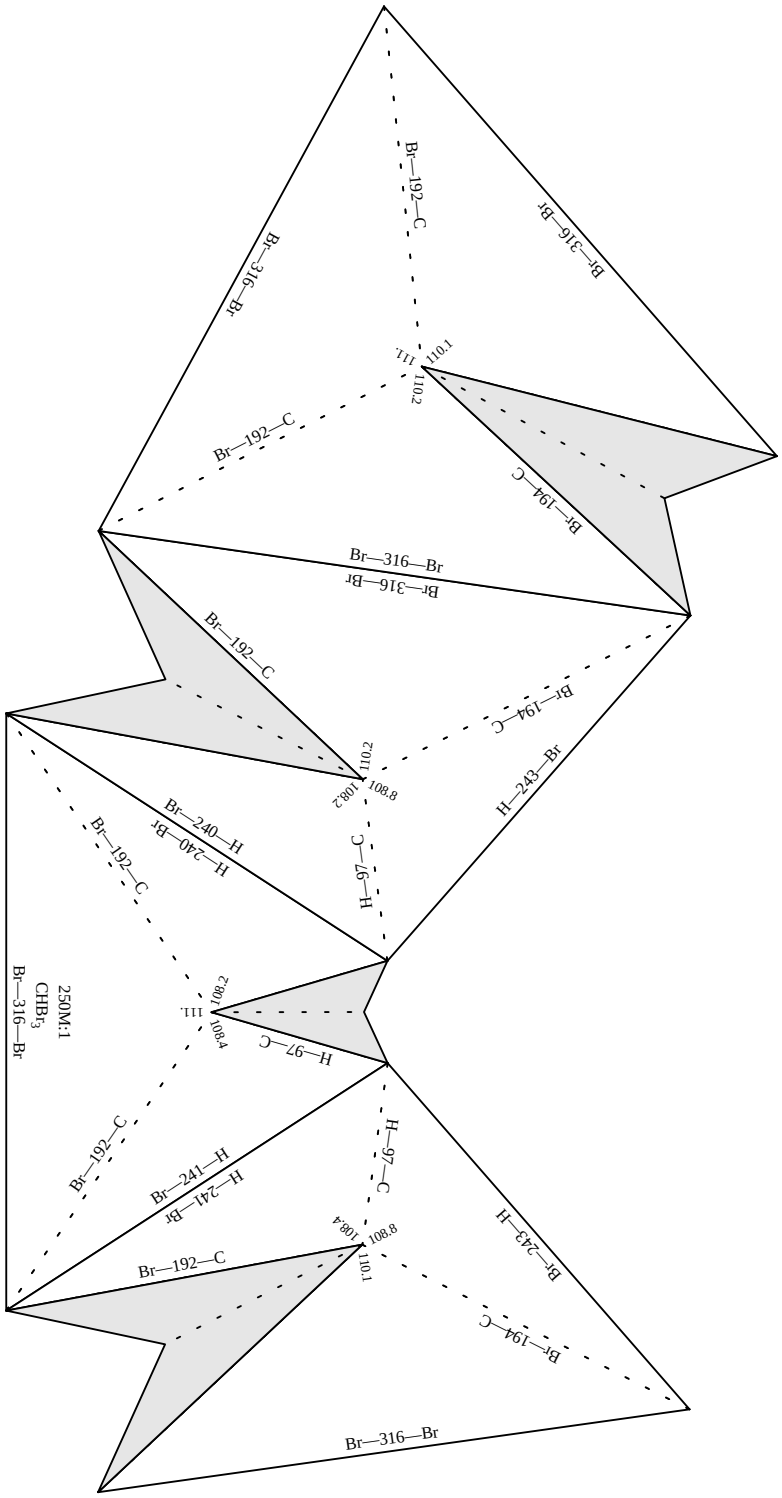
Molecular Origami of CHBr3

!C1
Br1
Br2
Br3
H1
CHBr3

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.99 offsety -0.47

View -1



Current: (centerx 3.31) (centery 4.53) (scale 250)
%%BoundingBox: 151 136 460 713 actual: 161 146 450 703 center: 306 425 actual size: 289 557
Better: (centerx 3.31) (centery 4.13) (scale 250)
%%BoundingBox: 80 74 389 651 actual: 90 84 379 641 center: 235 362 actual size: 289 557