Molecular Origami of CBr4

given information

ElementNames	[(C) (Br) (Br) (Br) (Br)]	
distance	185.595	C^1-Br^4
distance	185.595	C^1-Br^3
distance	185.595	C^1-Br^2
distance	224.786	C^1-Br^1
angle	100.990	Br^4-C^1-Br^1
	317.6	Br^4-Br^1
angle	100.990	Br^3-C^1-Br^1
	317.6	Br^3-Br^1
angle	100.990	Br^2-C^1-Br^1
	317.6	Br^2-Br^1
angle	116.455	Br^4-C^1-Br^3
	315.6	Br^4-Br^3
angle	116.455	Br^4-C^1-Br^2
	315.6	Br^4-Br^2
angle	116.455	Br^3-C^1-Br^2
	315.6	Br^3-Br^2
dopage	T	
AutoAlign	F	

structure type: XABCD

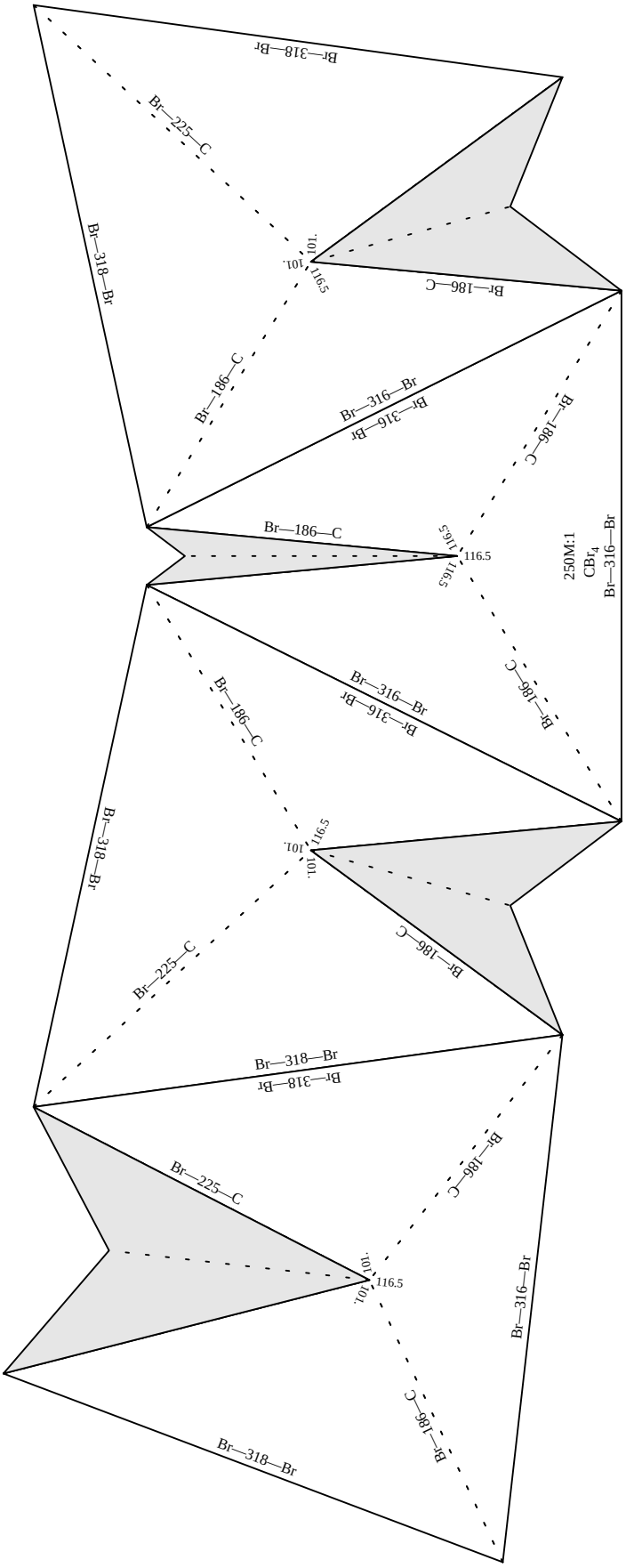
Molecular Origami of CBr4

!C1
Br1
Br2
Br3
Br4
CBr4

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx 0.8 offsety 2.23

View -1



Current: (centerx 5.10) (centery 7.23) (scale 250)
%%BoundingBox: 166 87 446 763 actual: 176 97 436 753 center: 306 425 actual size: 261 656
Better: (centerx 5.10) (centery 6.83) (scale 250)
%%BoundingBox: 223 218 504 895 actual: 233 228 494 885 center: 364 557 actual size: 261 656