Molecular Origami of AsBr3

given information

ElementNames	[(As) (Br) (Br) (Br)]	
dotted	F	
distance	232.175	As^{1}-Br^{3}
distance	232.179	As^{1}-Br^{1}
distance	232.182	As^{1}-Br^{2}
angle	98.927	Br^{2}-As^{1}-Br^{1}
	352.9	Br^{2}-Br^{1}
angle	98.927	Br^{3}-As^{1}-Br^{1}
	352.9	Br^{3}-Br^{1}
angle	98.928	Br^{3}-As^{1}-Br^{2}
	352.9	Br^{3}-Br^{2}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

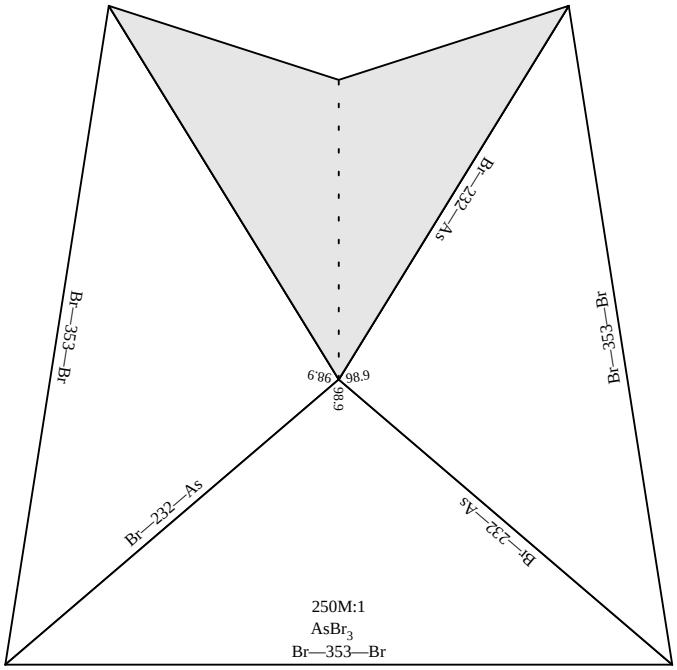
Molecular Origami of AsBr3

!As1
Br1
Br2
Br3
AsBr3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx 1.67 offsety 0.9

View -1



Current: (centerx 5.97) (centery 5.90) (scale 250)
%%BoundingBox: 170 290 440 557 actual: 180 300 430 547 center: 305 423 actual size: 250 247
Better: (centerx 5.99) (centery 5.52) (scale 250)
%%BoundingBox: 291 327 561 594 actual: 301 337 551 584 center: 426 461 actual size: 250 247