

Molecular Origami of SbPh2Br3

given information

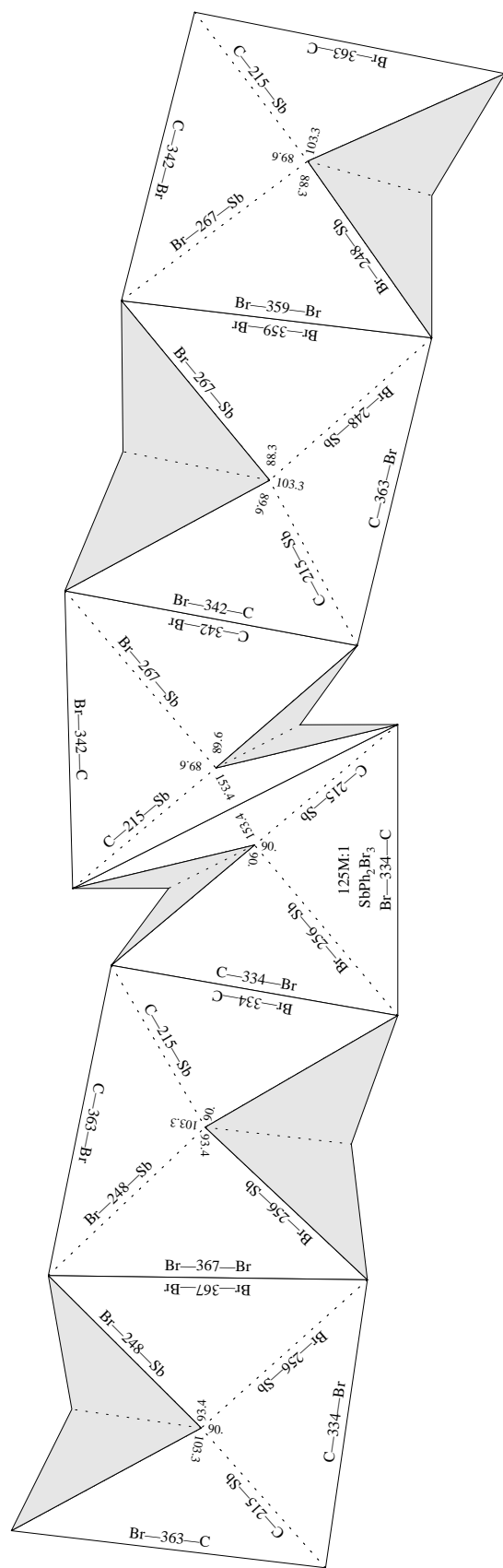
ElementNames	[(Sb) (Br) (Br) (Br) (C) (C)]	
distance	214.884	Sb ¹ -C ¹
distance	214.894	Sb ¹ -C ⁷
distance	247.899	Sb ¹ -Br ¹
distance	255.762	Sb ¹ -Br ²
distance	267.284	Sb ¹ -Br ³
angle	88.291	Br ³ -Sb ¹ -Br ¹
	359.1	Br ³ -Br ¹
angle	89.593	C ¹ -Sb ¹ -Br ³
	341.8	C ¹ -Br ³
angle	89.593	C ⁷ -Sb ¹ -Br ³
	341.8	C ⁷ -Br ³
angle	90.020	C ⁷ -Sb ¹ -Br ²
	334.1	C ⁷ -Br ²
angle	90.020	C ¹ -Sb ¹ -Br ²
	334.1	C ¹ -Br ²
angle	93.391	Br ² -Sb ¹ -Br ¹
	366.6	Br ² -Br ¹
angle	103.287	C ⁷ -Sb ¹ -Br ¹
	363.5	C ⁷ -Br ¹
angle	103.288	C ¹ -Sb ¹ -Br ¹
	363.5	C ¹ -Br ¹
angle	153.376	C ⁷ -Sb ¹ -C ¹
	418.2	C ⁷ -C ¹
angle	178.318	Br ³ -Sb ¹ -Br ²
	523.	Br ³ -Br ²
dopage	T	
AutoAlign	F	

structure type: XABCDE

!Sb1
Br1
Br2
Br3
C1
C7
SbPh2Br3

```
scale 125,000,000 : 1
units: pm
offsetx -0.22 offsety 0.46
```

View -1



actual size: 201 633

actual size: 201 633