

Molecular Origami of Sb(CH3)4⁺

given information

ElementNames	[(Sb) (C) (C) (C) (C)]	
distance	207.947	Sb ¹ -C ³
distance	208.947	Sb ¹ -C ¹
distance	210.428	Sb ¹ -C ²
distance	213.662	Sb ¹ -C ⁴
angle	107.380	C ³ -Sb ¹ -C ¹
	335.9	C ³ -C ¹
angle	107.776	C ⁴ -Sb ¹ -C ¹
	341.4	C ⁴ -C ¹
angle	109.017	C ³ -Sb ¹ -C ²
	340.6	C ³ -C ²
angle	109.583	C ⁴ -Sb ¹ -C ³
	344.5	C ⁴ -C ³
angle	110.964	C ⁴ -Sb ¹ -C ²
	349.4	C ⁴ -C ²
angle	112.034	C ² -Sb ¹ -C ¹
	347.7	C ² -C ¹
dopage	T	
AutoAlign	F	

structure type: XABCD

!Sb1
C1
C2
C3
C4
Sb(CH3)4^+

```
scale 250,000,000 : 1
units: pm
offsetx -0.55 offsety -0.34
```

actual size: 236 740

actual size: 213 693