

Molecular Origami of N(CH3)4⁺

given information

ElementNames	[(N) (C) (C) (C) (C)]	
distance	147.934	N ¹ -C ⁴
distance	147.934	N ¹ -C ³
distance	147.967	N ¹ -C ¹
distance	149.465	N ¹ -C ²
angle	106.535	C ⁴ -N ¹ -C ³
	237.1	C ⁴ -C ³
angle	109.819	C ⁴ -N ¹ -C ¹
	242.1	C ⁴ -C ¹
angle	109.819	C ³ -N ¹ -C ¹
	242.1	C ³ -C ¹
angle	109.847	C ⁴ -N ¹ -C ²
	243.4	C ⁴ -C ²
angle	109.847	C ³ -N ¹ -C ²
	243.4	C ³ -C ²
angle	110.877	C ² -N ¹ -C ¹
	244.9	C ² -C ¹
dopage	T	
AutoAlign	F	

structure type: XABCD

!N1
C1
C2
C3
C4
N(CH3)4^+

```
scale 250,000,000 : 1
units: pm
offsetx -0.33 offsety 0.05
```

actual size: 167 518

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