Molecular Origami

given information

| | | |
|---|---|---|
| ElementNames | [ (P) (C) (C) (C) ] | |
| dotted | F | |
| distance | 183.297 | $P^1$-$C^1$ |
| distance | 183.669 | $P^1$-$C^{10}$ |
| distance | 183.796 | $P^1$-$C^{19}$ |
| angle | 107.921 | $C^{19}$-$P^1$-$C^1$ |
| | 296.8 | $C^{19}$-$C^1$ |
| angle | 110.126 | $C^{10}$-$P^1$-$C^1$ |
| | 300.8 | $C^{10}$-$C^1$ |
| angle | 110.666 | $C^{19}$-$P^1$-$C^{10}$ |
| | 302.2 | $C^{19}$-$C^{10}$ |
| dopage | T | |
| AutoAlign | F | |
| showboth | T | |

structure type: XABC

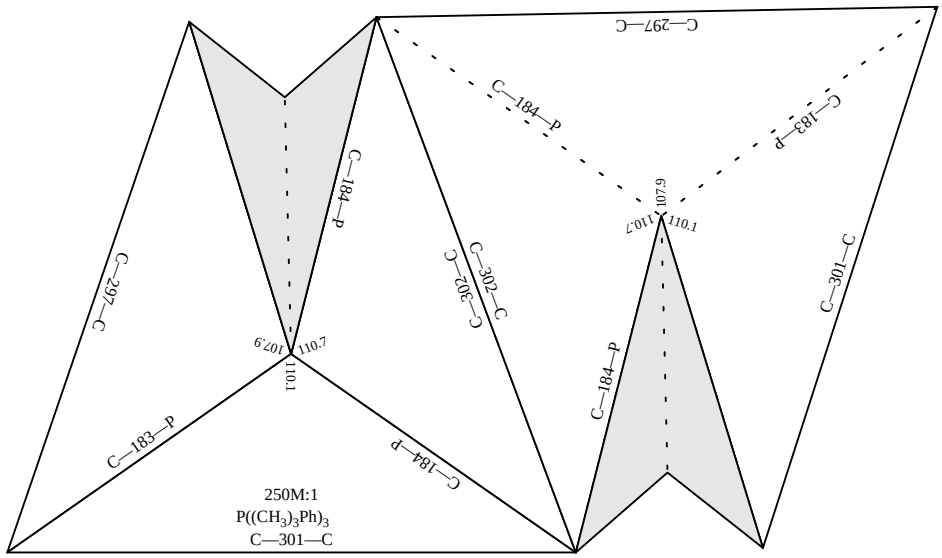
Molecular Origami

!P1
C1
C10
C19
P((CH3)3Ph)3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx 1.35  offsety 0.9

View -1




Current:  (centerx 5.65) (centery 5.90) (scale 250)
%%BoundingBox: 184 308 552 533  actual: 194 318 542 523  center: 368 421  actual size: 349 205
Better:  (centerx 4.79) (centery 5.56) (scale 250)
%%BoundingBox: 219 348 588 573  actual: 229 358 578 563  center: 403 461  actual size: 349 205