

Supporting Information

New fluorescence labeling isotactic polypropylenes as a tracer: a proof of concept

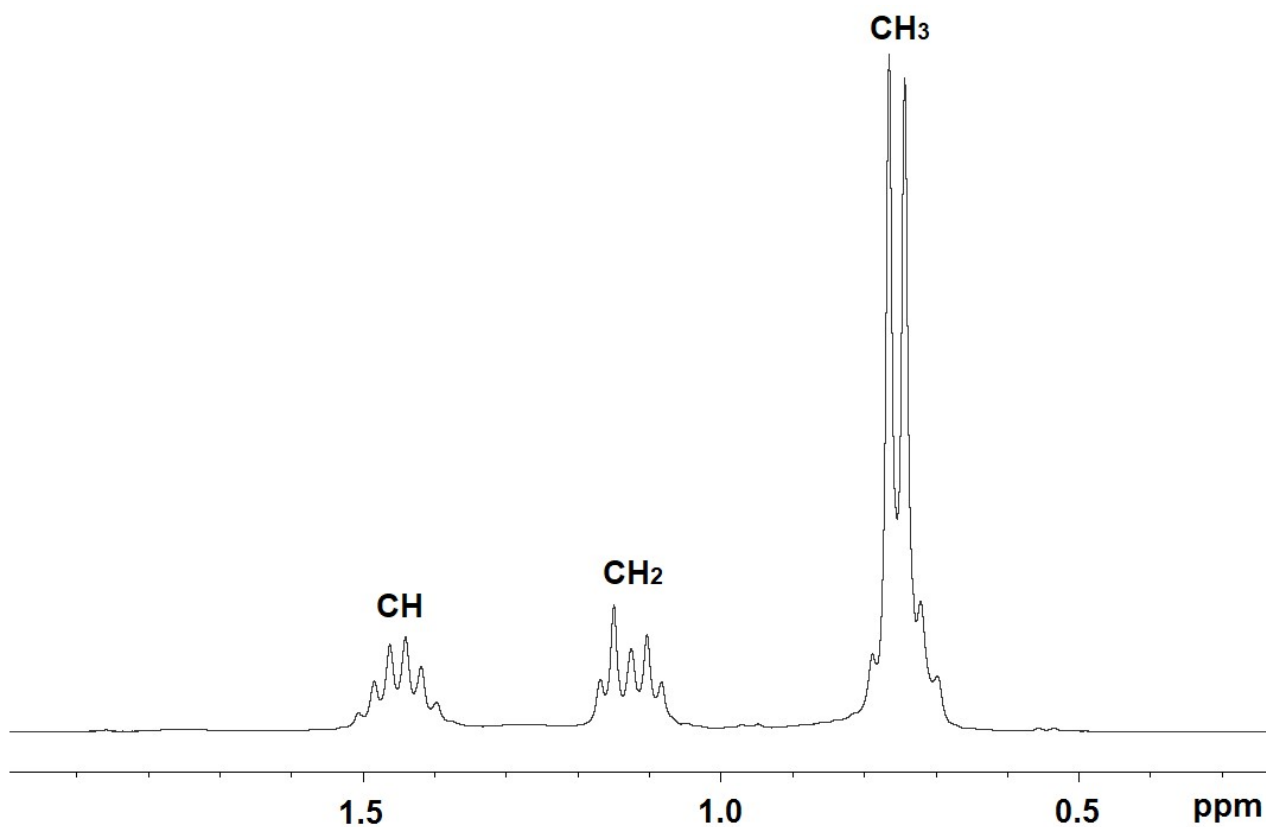
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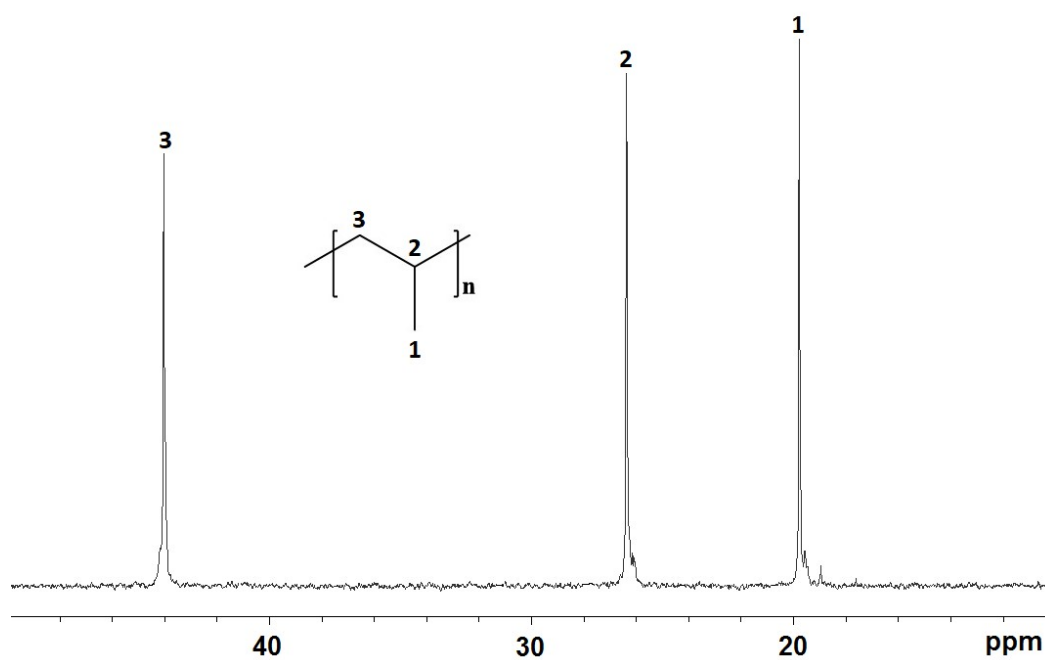
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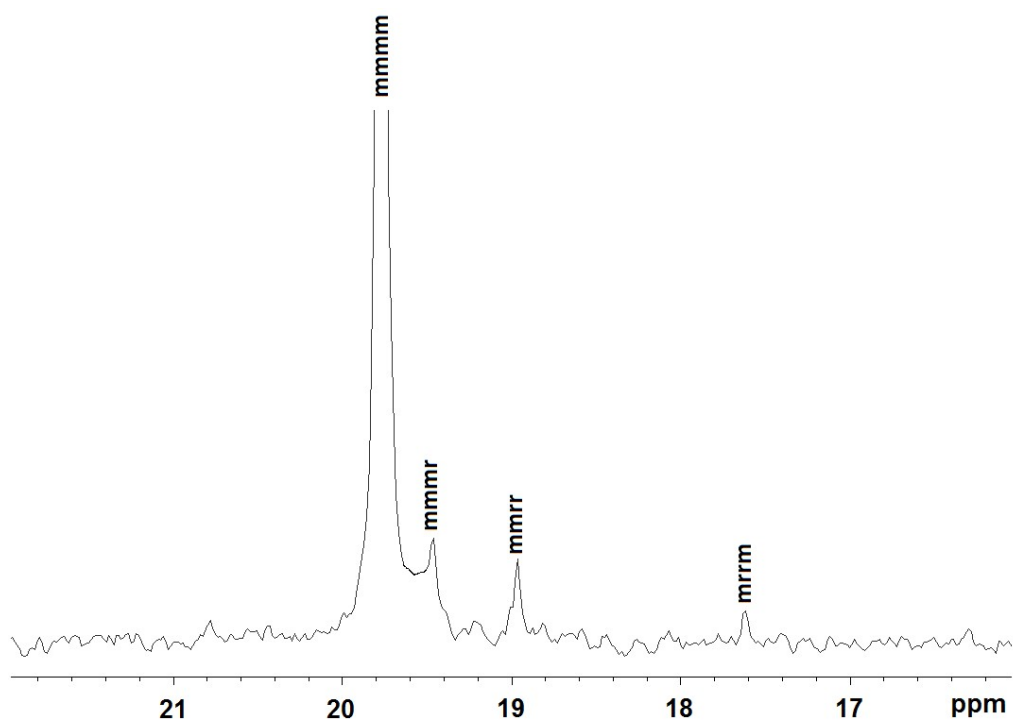
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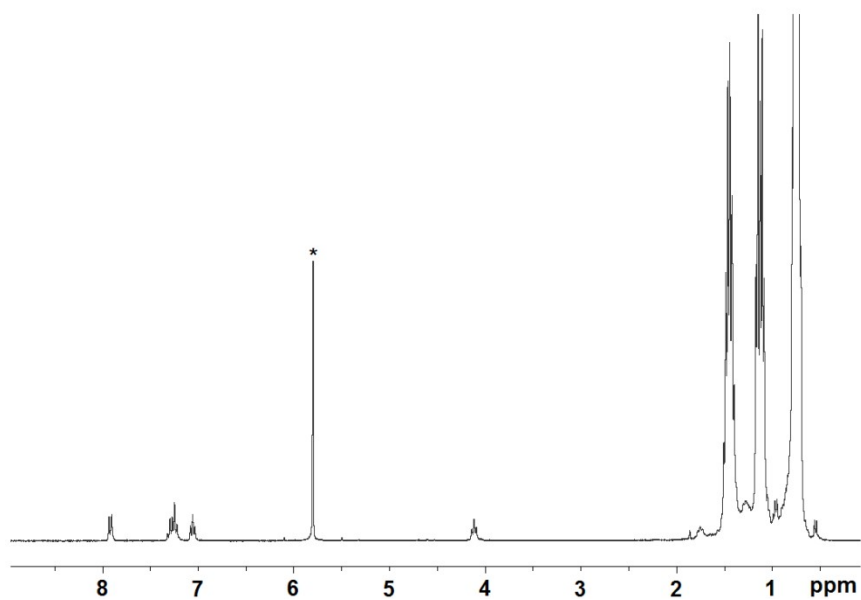
S1. ^1H NMR spectrum of *i*-PP (run 1) (TCDE solvent, HMDS scale, 100 °C).



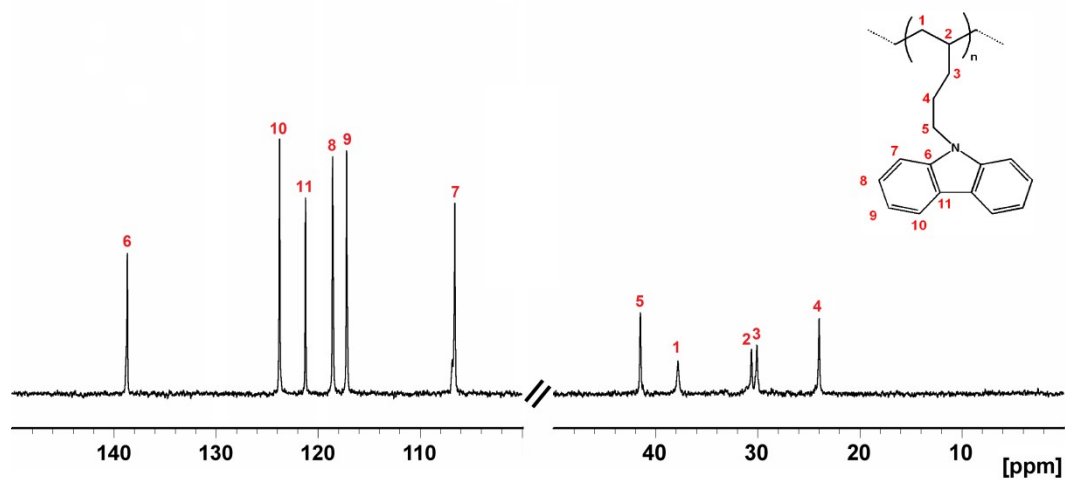
S2. ^{13}C NMR spectrum of *i*-PP (run 1) (TCDE solvent, HMDS scale, 100 °C).



S3. Methyl region of ^{13}C NMR spectrum of *i*-PP (run 1) (TCDE solvent, HMDS scale, 100 °C).



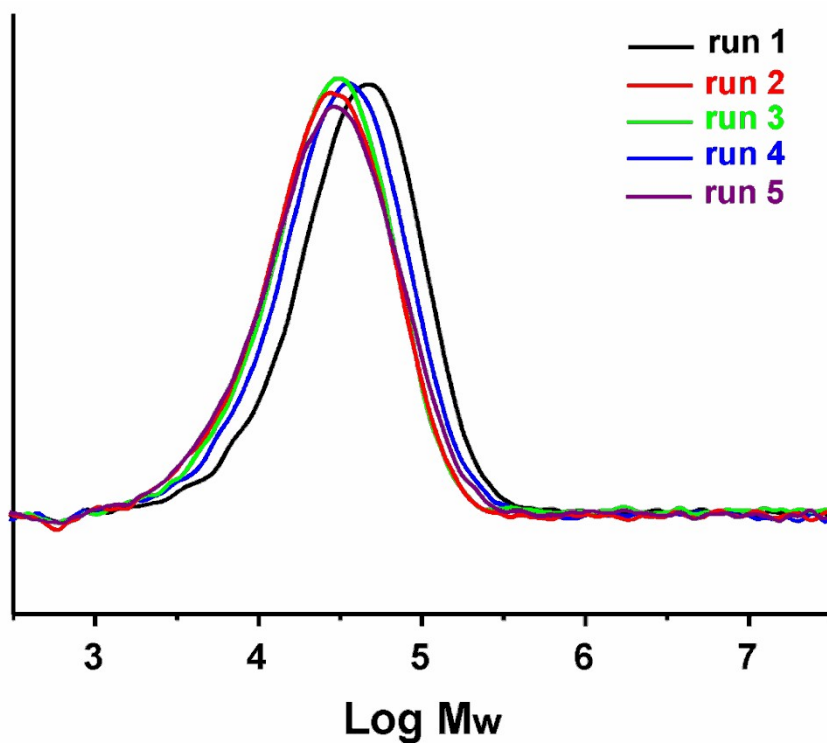
S4. ^1H NMR spectrum of P(P-co-PK) (run 2) (TCDE solvent, HMDS scale, 100 °C).



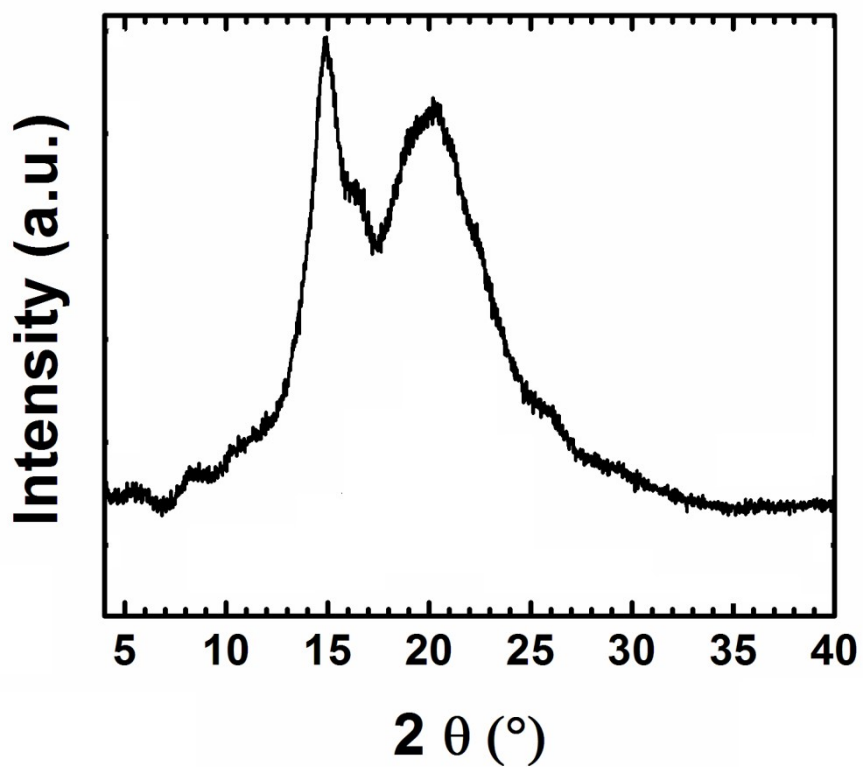
S5. ¹³C NMR spectrum of PPK (run 5) (TCDE solvent, HMDS scale, 100 °C).

Table S1. ¹³C NMR Assignments of Run 5

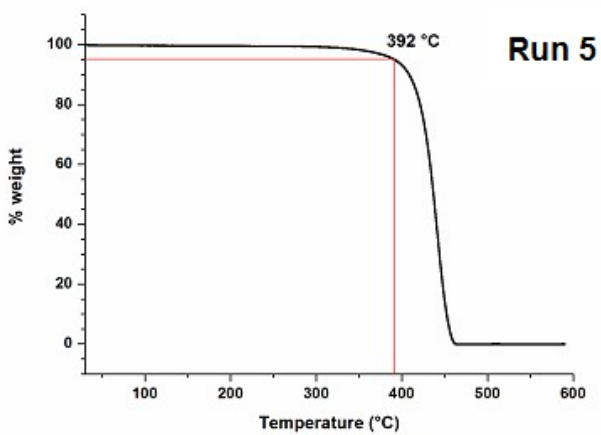
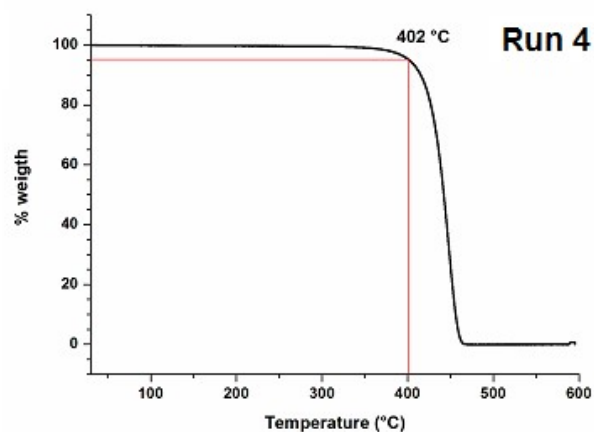
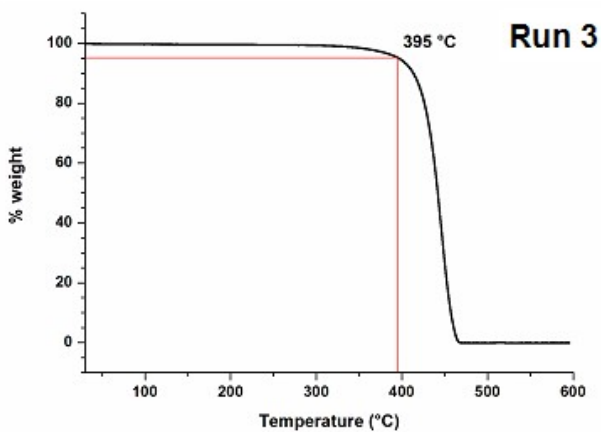
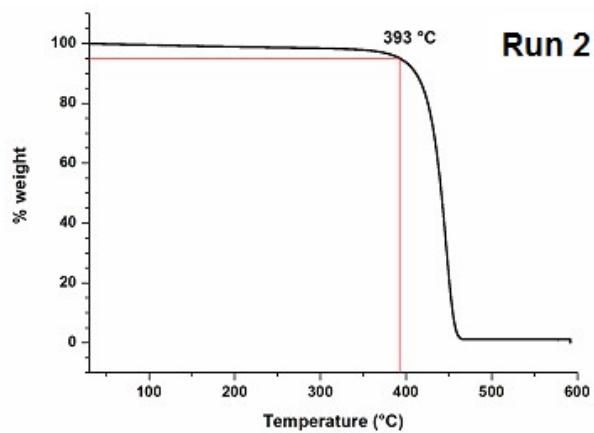
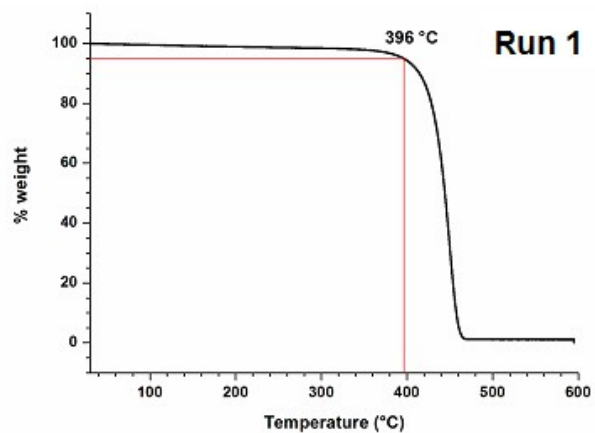
Carbon	Chem. Shift (ppm)
C1	37.72 ₆
C2	30.53 ₈
C3	23.94 ₆
C4	30.02 ₂
C5	41.45 ₆
C6	138.63 ₉
C7	106.57 ₁
C8	118.55 ₆
C9	117.17 ₇
C10	123.75 ₆
C11	121.17 ₁



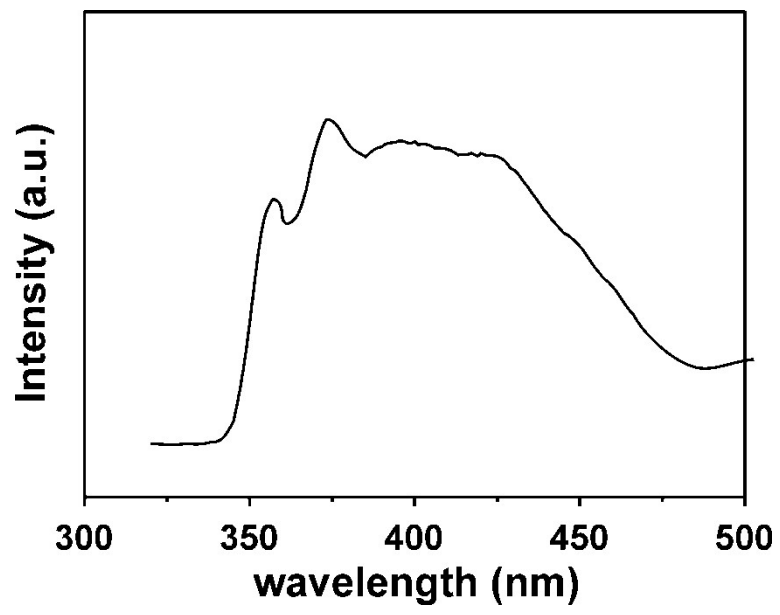
S6. GPC traces of runs 1-5.



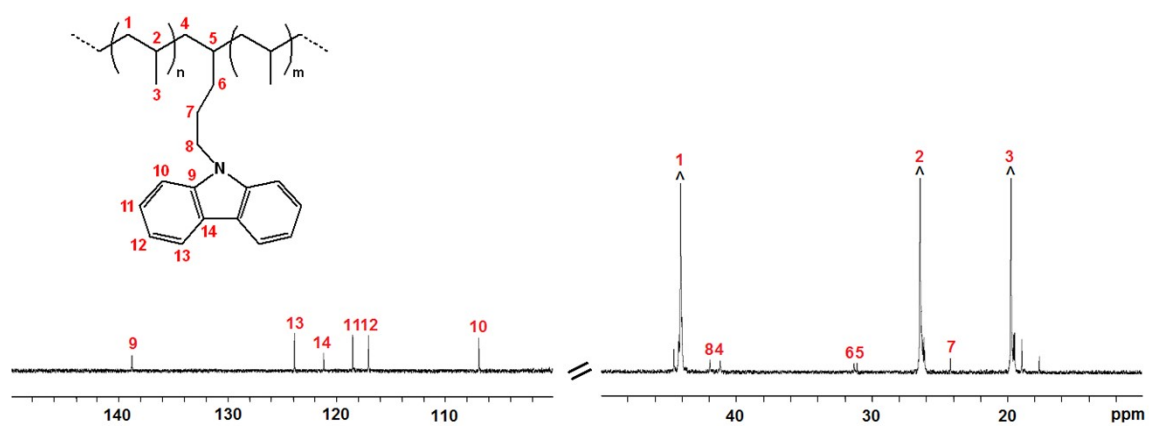
S7. X ray diffraction pattern of PPK (run 5).



S8. TGA traces of runs 1-5 (N₂, 10°C/min).



S9. Emission spectrum of PPK (run 5). $\Lambda_{\text{ecc}} = 300$ nm.



^{13}C NMR spectrum of P(P-co-PK) (run 2) (TCDE solvent, HMDS scale, 100 °C).

Table S2. ^{13}C NMR Assignments of Run 2

Carbon	Chem. Shift (ppm)
C1	44.06 ₈
C2	26.41 ₄
C3	19.79 ₉
C4	41.11 ₉
C5	31.01 ₇
C6	31.28 ₆
C7	24.26 ₉
C8	41.96 ₇
C9	138.77 ₁
C10	106.93 ₁
C11	118.52 ₆
C12	117.05 ₆
C13	123.87 ₁
C14	121.13 ₂