Electronic Supplementary Material (ESI) for Chemical Communications.

Supporting Information

SUPPORTING INFORMATION

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SUPPORTING INFORMATION

Experimental Procedures

General Methods

NMR spectra were recorded on a Bruker Advance III 400 MHz, a Bruker Advance HD 400 and an Agilent 400 spectrometer for ^{1}H (400.13 MHz) and ^{13}C (100.61 MHz) NMR spectra. A Bruker Ultrashield 600 spectrometer was employed for ^{1}H (600.13 MHz), ^{13}C (150.90 MHz) and ^{15}N NMR (61 MHz) spectra. All NMR experiments were performed at 25 ºC. Resonances δ are given in ppm units and referenced to the deuterium signal in the NMR solvents, acetonitrile-d_3 (δ_H = 1.94 ppm, δ_C = 1.32, 118.26 ppm). Signal multiplicities are abbreviated as follows: singlet = s, doublet = d, triplet = t, dq = doublet of quartets, multiplet = m.

Normal-structural decomposition (NSD): The NSD method, as developed by Shelnutt and coworkers,[1] was used to delineate, quantify, and illustrate the various distortions modes present in the tetrapyrrole macrocycles. Analysis was performed with the NSD online interface, available at https://www.sengegroup.eu/nsd.[2]

Single crystal X-ray crystallography: Crystals were grown following the protocol developed by Hope, liquid-liquid diffusion of $CHCl_3$ and MeOH with H_2SO_4.[3] Diffraction data were collected on a Bruker APEX 2 DUO CCD diffractometer using Incoatec IμS Cu-K_α (λ = 1.54178 Å) radiation. Crystal was mounted on a MiTeGen MicroMount and collected at 100(2) K using an Oxford Cryosystems Cobra low-temperature device. Data were collected using omega and phi scans and were corrected for Lorentz and polarization effects using the APEX software suite.[4] Data were corrected for absorption effects using the multi-scan method (SADABS).[5] Using Olex2, the structure was solved with the XT structure solution program, using the intrinsic phasing solution method and refined against $|F^2|$ with XL using least-squares minimization.[6] If electron density was not sufficient, the C and N bound H atoms were placed in their expected calculated positions and refined using a riding model: N–H = 0.88 Å, C–H = 0.95–0.98 Å, with U_{iso} (H) = 1.5U_{eq} (C) for methyl H atoms and 1.2U_{eq} (C, N). Details of data refinements can be found in Table S4. All images were prepared using Olex2.[6a]

In the structure $\alpha_2\beta_2$-**P**[SO_4^{2-}][HSO_4^-]$_4$, two phenyl rings at C5 and C10 and one ethyl group are modelled over two locations using DFIX, SIMU, SADI restrains and EADP constraints. In terms of counter anion, only the SO_4^{2-} group is not disordered, while all other HSO_4^- groups were modelled disordered using rigid groups. Some hydrogen atoms were placed geometrically to compensate for close contacts, the remaining hydrogens could not be located on the disordered HSO_4^- moieties but were added to the formula to make the formula weight correct. Multiple disordered and partially occupied H_2O molecules are modelled in the structure using SIMU and ISOR restrains. The weighting scheme was manually adjusted to ensure the goodness of fit was reasonable.

Deposition number 2143572 contain the supplementary crystallographic data for this paper. These data are provided free of charge by the joint Cambridge Crystallographic Data Centre and Fachinformationszentrum Karlsruhe http://www.ccdc.cam.ac.uk/structures.

All calculations were done with Gaussian16 Rev. B.01.[7] A conformational search for $\alpha_2\beta_2$-**P·10CSA(R)** was performed using a crystal structure of $\alpha_2\beta_2$-**P**[SO_4^{2-}][HSO_4^-]$_4$ supramolecules as a starting point for building porphyrin/camphorsulfonic acid supramolecules, followed by optimization in acetonitrile. The geometry optimization and frequencies calculations were performed using BP86[8]-D3BJ[9]/def2-SVP[10] – the method which showed a good agreement with the experimental data reported in our previous works.[11] To include acetonitrile effects the SMD[12] continuum solvent model was used, where molecular surface was represented as Solvent Accessible Surface (SAS) and the Bondi atomic radii were used. A ground state was characterized by absence of imaginary frequencies and more accurate electronic energies were calculated using the BP86-D3BJ/def2-TZVPP[10] and SMD model. During conformational search twenty start geometries converged into eight conformers corresponding to the ground state (Table S5) with one major conformer A making up 90%. The geometry were visualized using GaussView 6.1.[13] and is illustrated in figure S14.

The NMR shielding tensors were calculated at the GIAO[14]-B3LYP[15]/6-311++G** [14a, 16] level of theory using the SMD continuum solvent model to include acetonitrile effects. To calculate the ^{13}C chemical shifts, a scaling factor of 1.0228[17] and a reference point TMS (180.7 ppm) were used , for calculation of the ^{15}N chemical shifts NH_3 was used as a reference point (253.70 ppm) (Table S6 – S8). Population analysis was done using the BP86-D3BJ/def2-TZVPP//BP86-D3BJ/def2-SVP level of theory in acetonitrile and NBO[18] approach in acetonitrile. Non-covalent interactions were also analyzed using the SMD (acetonitrile), the BP86-D3BJ/def2-SVP level of theory and AIMAll program version 19.10.12[19]

In Text Supplementary Material

*NMR Investigation of the **P-10CSA** Atropisomers*

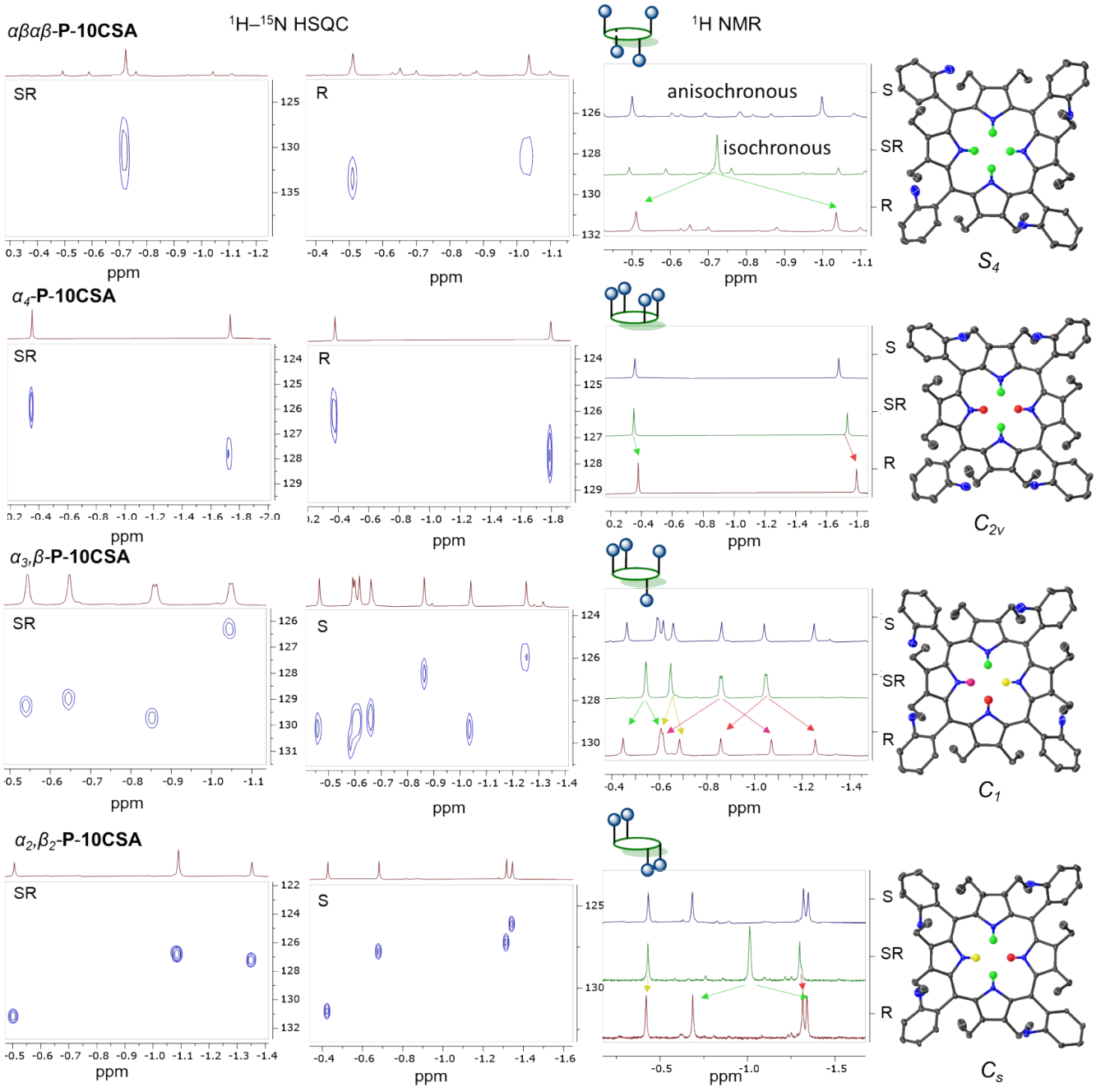


Figure S1. ^{1}H and ^{1}H–^{15}N HSQC NMR spectra obtained for 0% and 100% e.e. **P·10CSA** solutions (20 equivalents) in d_3-acetonitrile. On the right side, a graphical illustration of **P·10CSA** atropisomers with corresponding point group notations, inner core system protons are highlighted in different colors correlating to the arrows marked in ^{1}H NMR spectra.

We performed ^{1}H NMR and ^{1}H–^{15}N HSQC analyses of all **P** atropisomers with racemic mixtures and enantiopure **10CSA**(**S** or **R**). For the racemic **10CSA** solutions, as expected ^{1}H NMR spectra of **P** atropisomers remained isochronous (Figure S1). Due to the chiral information transmitted in equal proportions from both chiral components (**10CSA R** and **S**), the observed inner core signals resonate in an identical manner to achiral acids (**BSA** and **MSA**) previously reported by us.[20] One singlet is observed for *αβαβ*-**P·10CSA**, two in α_4-**P·10CSA**, three signals of relative intensity 1:2:1 in $\alpha_2\beta_2$-**P·10CSA**, and finally the spectrum for the unsymmetrical $\alpha_3\beta$-**P·10CSA** atropisomer has four differently shifted signals. The enantiopure solutions are described in the main text.

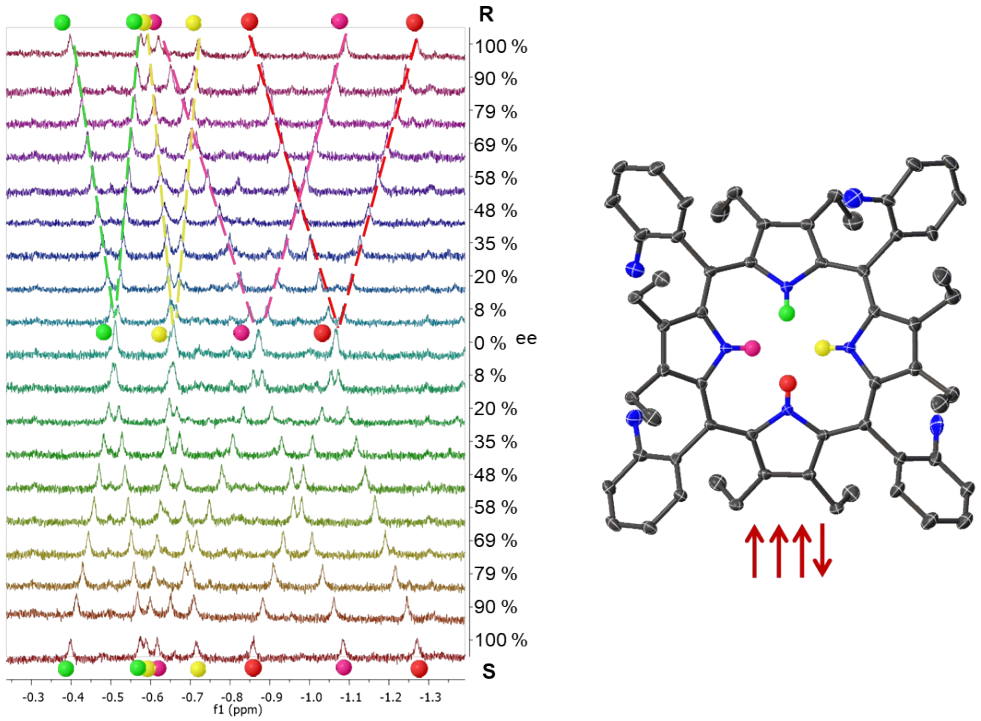


Figure S2. ^{1}H NMR titration studies of the $\alpha_3\beta$-**P** inner core system using different ratios of chiral components (**10CSA R** and **S**, 20 eq.). Spectra recorded in CD_3CN

10CSA Additions to $\alpha_2\beta_2$-*P*

The gradual addition of **10CSA(R)** into the CD_3CN solution of $\alpha_2\beta_2$-**P** resulted in downfield shifts of the ^{1}H NMR aromatic region signals with the and observable splitting pattern of the *o*-Ar*H* (Figure S3), while the split pyrrolic N–*H* were upfield shifted upon emerging at ~3 eq. of **10CSA(R)** in the ^{1}H NMR spectra (Figures S4 and S5). The magnitude of chemical shifts of split peaks (Δσ) drastically increased over the first ~7 eq., while the later additions resulted in marginal changes to Δσ (Figure S6).

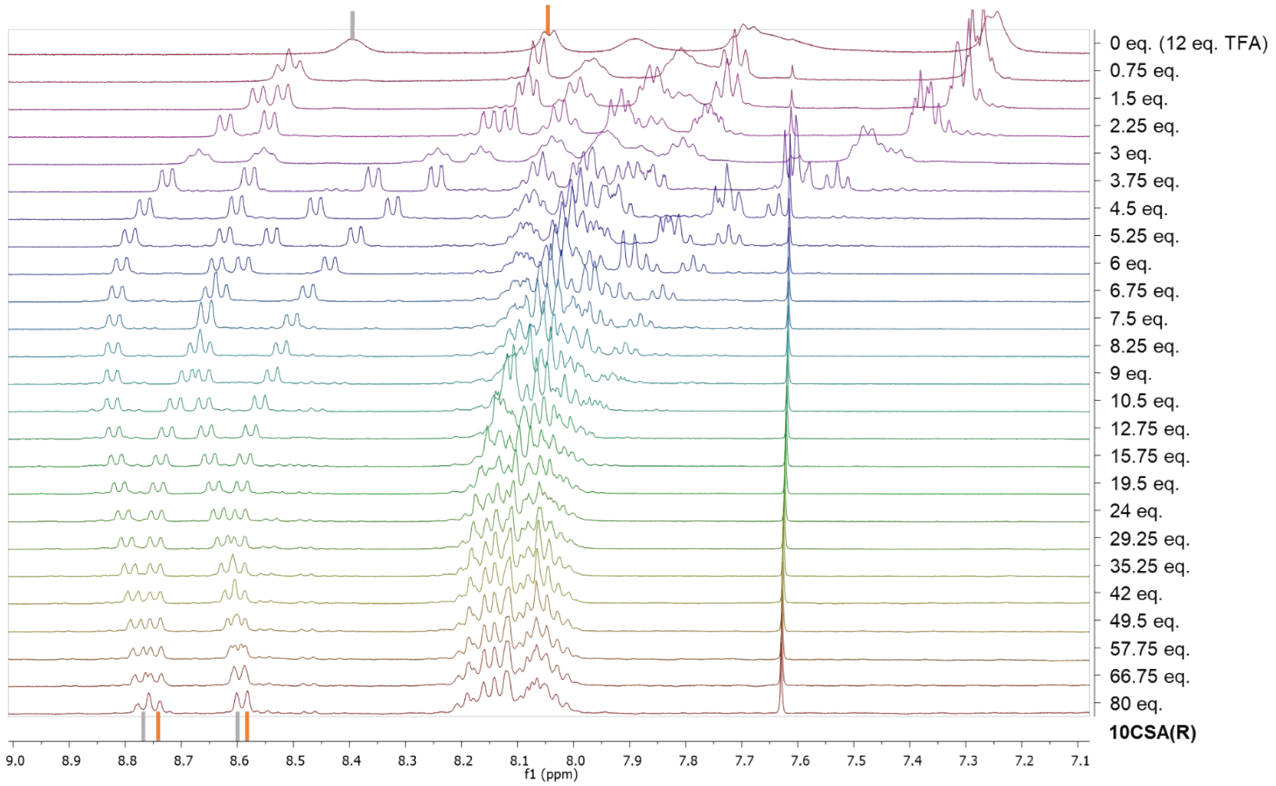


Figure S3. ^{1}H NMR of the $\alpha_2\beta_2$-**P** aromatic region dependence to the **10CSA(R)** equivalents, recorded in CD_3CN. The addition of 12 eq. of TFA was to solubilize and protonate $\alpha_2\beta_2$-**P**. Highlighted in grey and orange is two different splitting *o*-Ar-*H* regions.

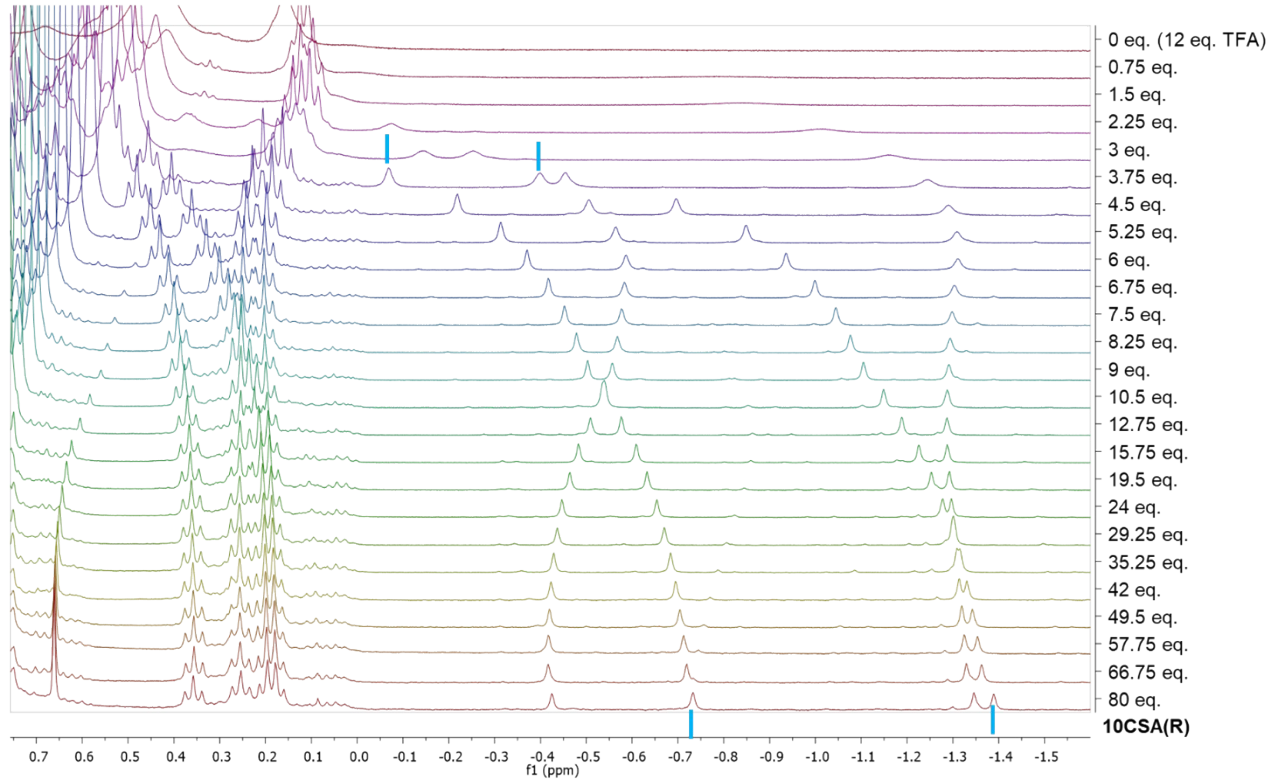


Figure S4. ^{1}H NMR of the $\alpha_2\beta_2$-**P** N–*H* and CH_3 regions dependence to the **10CSA(R)** equivalents, recorded in CD_3CN. The addition of 12 eq. of TFA was to solubilize and protonate $\alpha_2\beta_2$-**P**. Highlighted in blue are two different splitting N–*H* resonance signals.

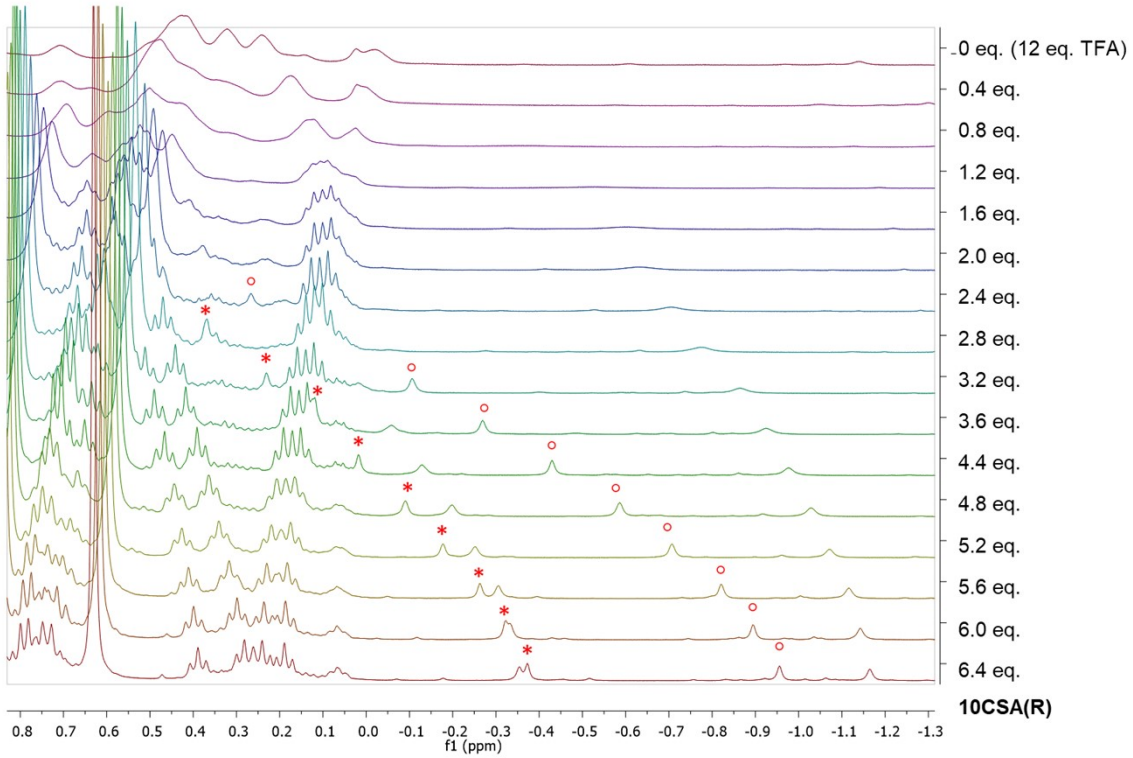


Figure S5. ^{1}H NMR of the $\alpha_2\beta_2$-**P** N–*H* regions dependence to the **10CSA(R)** equivalents, recorded in CD_3CN. The addition of 12 eq. of TFA was to solubilize and protonate $\alpha_2\beta_2$-**P**. Highlighted with red dot and star is two different splitting N–*H* resonance signals. Note, these spectra were recorded with higher concentrations of $\alpha_2\beta_2$-**P** expecting to get a better resolution of the inner N–*H*. From the obtained spectra it appears that the inner core system of interest originates downfield shifted and overlays with the aliphatic C*H*$_3$ region.

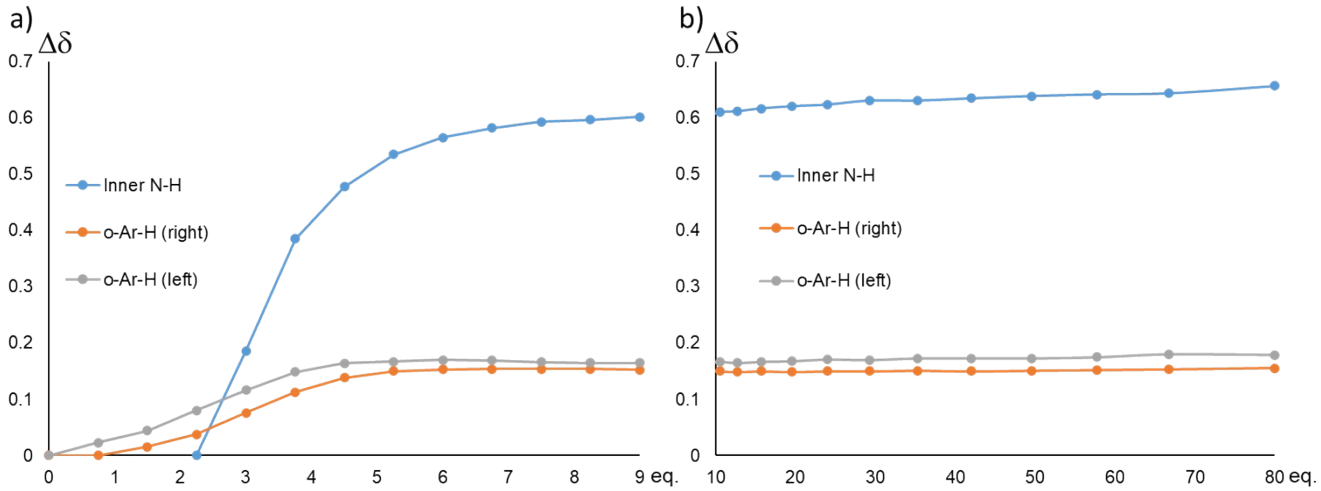


Figure S6. A graph showing the magnitude of the split chemical shifts Δδ in the inner N–*H* and *o*-Ar-*H* of the phenyl rings independence on the **10CSA(R)** equivalents, a) 0 to 9 eq., b) 10 to 80 eq. Recorded in CD_3CN with the addition of 12 eq. of TFA to solubilize and protonate $\alpha_2\beta_2$-**P**. The color coding corresponds to the highlighted signals in Figures S3 and S4.

Influence of Water on the Complexation

We have previously mentioned that one of the best-known *pro*-CSA (**Bz_2oxP**) suffers serious sensitivity issues due in part to *N*-alkylation.[21] Additionally, the competitive binding of water molecules significantly contributes to the magnitude of splitting Δσ. Trace amounts of water in solutions or titrants results in the necessary use of high guest concentrations to obtain well-resolved spectra. The hypersensitivity towards water is a major limitation for functional *pro*-CSA, since water is ubiquitous, avoiding it is at least tedious if not almost impossible for most of the solvents and reagents, especially for the analytes bearing high polarity. Hence, investigation of water influence as a competitive agent to the enantio-pure $\alpha_2\beta_2$-**P·10CSA(S)** complex in CD3CN was carried out. From the first instance, the gradual addition of water has a considerable effect on Δσ. By the addition of 240 eq. of water, the Δσ of N–H has contracted by 0.132 ppm. (Figure S9). Despite the competitive nature of water, the overall binding strength of **10CSA(S)** remains observable considering the ~12.5 fold higher water consistency in the solution. Even at substantially higher (>4000 eq.) quantity of water the signals remained anisochronous highlighting the strong relationship between the host and chiral guest (Figure S8-S9).

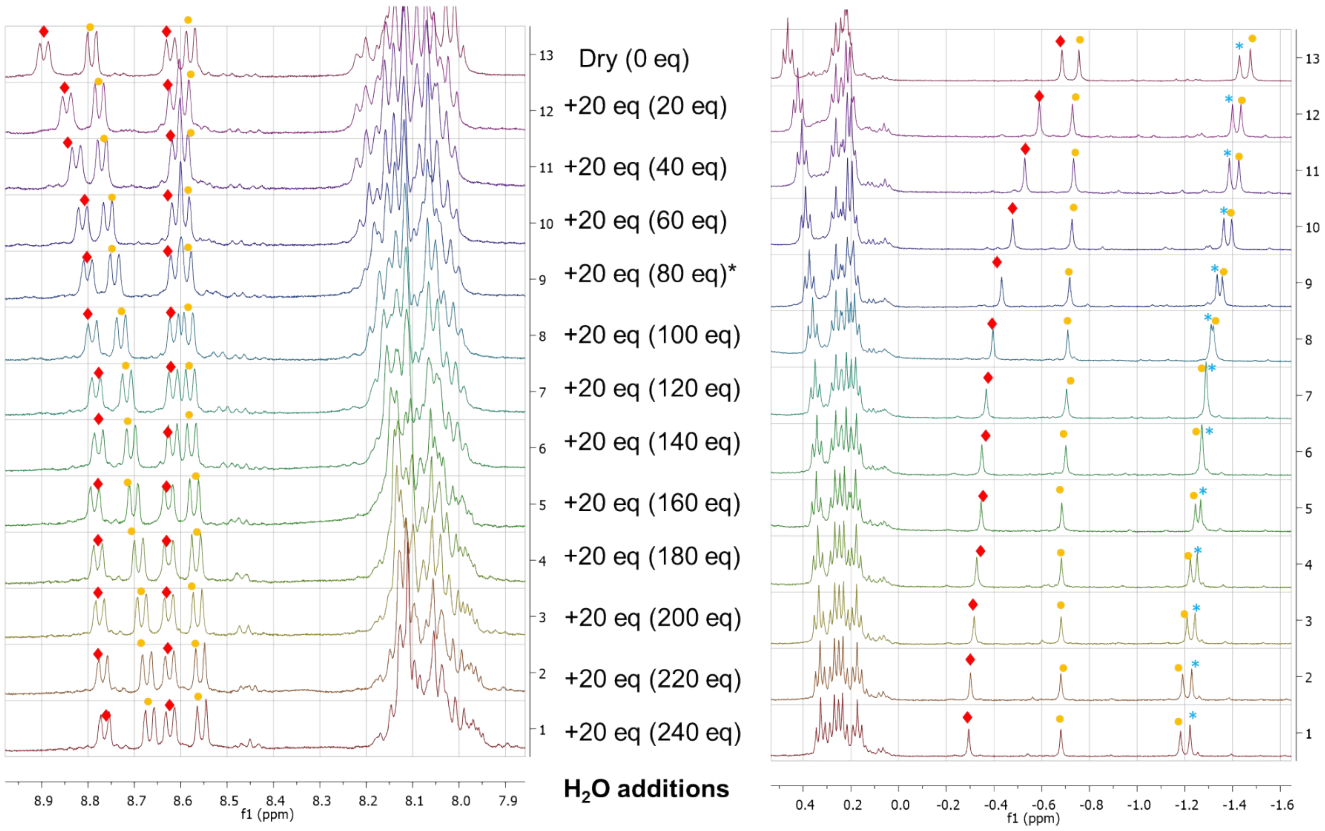


Figure S7. ^{1}H NMR titration with H_2O (0-240 eq) of $\alpha_2\beta_2$-**P·10CSA(S)** (19 eq). The split inner N–*H* is highlighted with orange dots. In the aromatic region, the split *o*-Ar*H* signals are highlighted with orange and red dots. Spectra recorded in CD_3CN

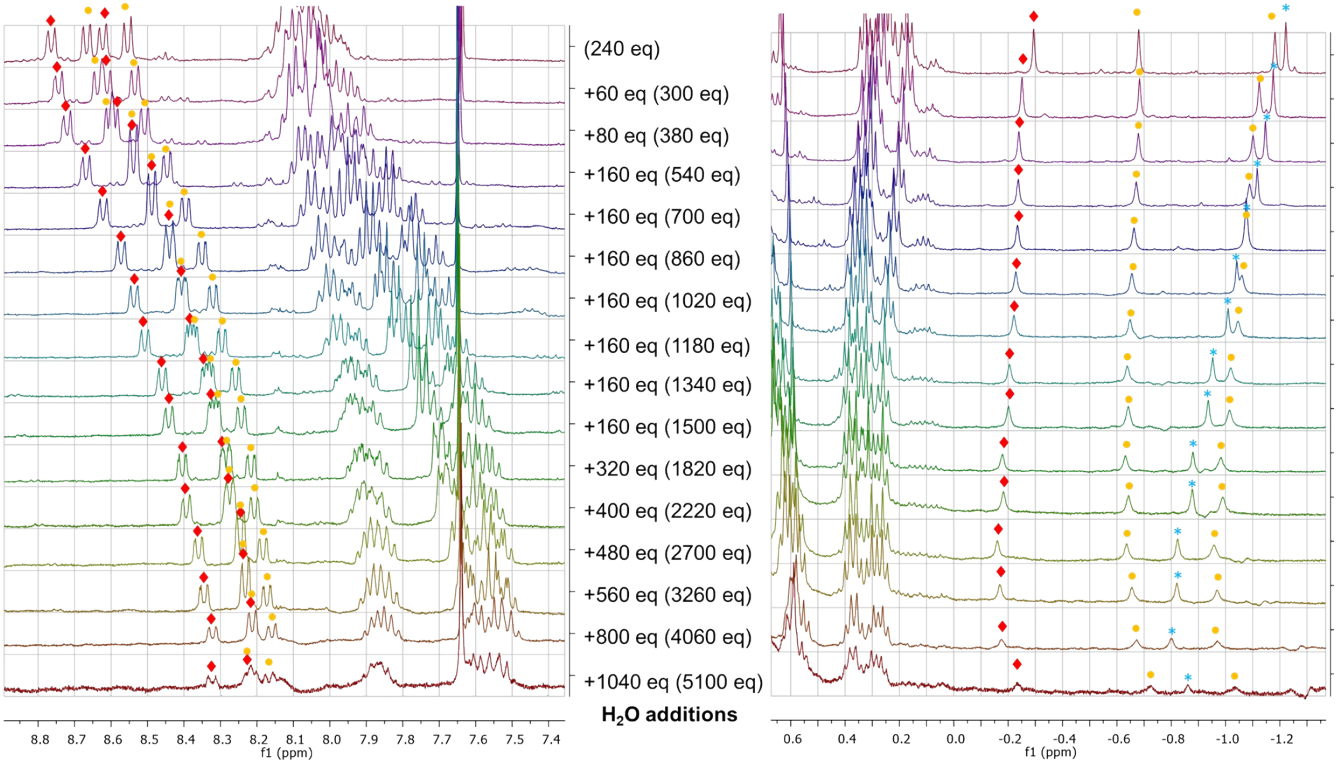


Figure S8. ^{1}H NMR titration with H_2O (240-5100 eq) of $\alpha_2\beta_2$-**P·10CSA(S)** (19 eq). The split inner N–*H* is highlighted with orange dots. In the aromatic region, the split *o*-Ar*H* signals are highlighted with orange and red dots. Spectra recorded in CD_3CN.

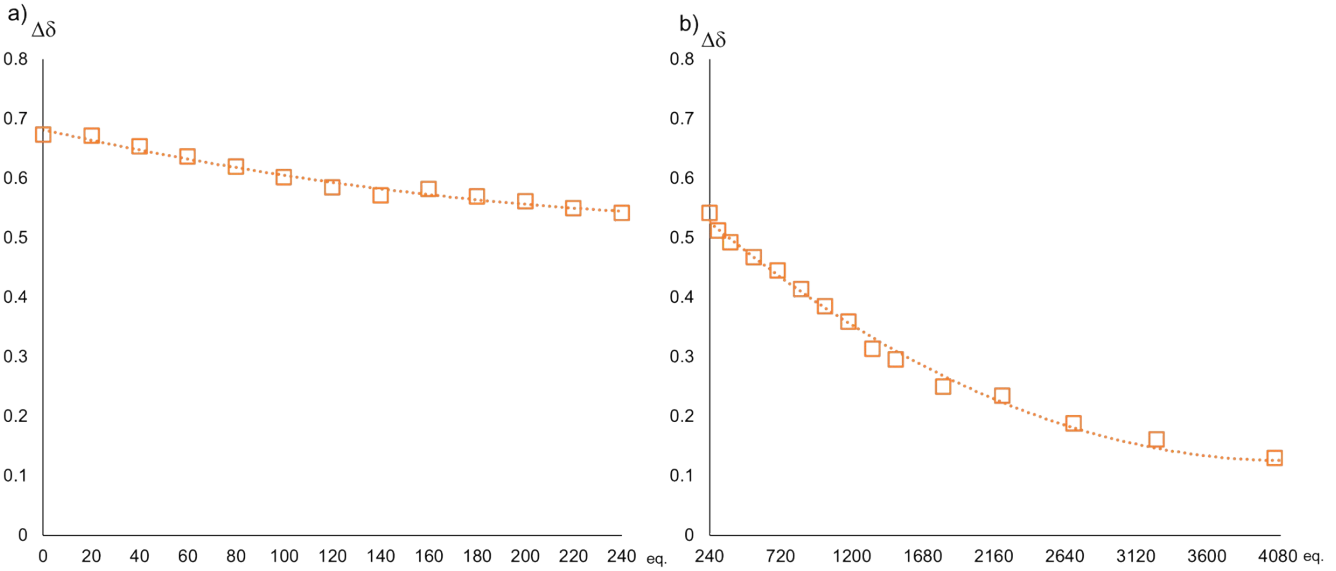


Figure S9. The graph of $\alpha_2\beta_2$-**P·10CSA(S)** (19 eq) representing N–*H* Δσ dependence to the various additions of water a) 0-240 eq. b) 240-5100 eq.

Detection of Enantiomeric Excess

The origin of the chemical shift non-equivalence lies deep within the concept of prochirality.[22] The desymmetrization of $\alpha_2\beta_2$-**P** atoms in a single step by weak interactions with a chiral guest proves to be particularly useful for the e.e. determination. When racemic and non-racemic mixtures of **10CSA** were applied in the system at constant concentrations it was found that the magnitude of the split $\Delta\sigma$ of the *o*-Ar*H*, C*H*$_3$, and N–H peaks rely on respective e.e. value (%) (Figure 2c). At the racemate point, the isochronous profile of $\alpha_2\beta_2$-**P** is restored since the chiral information is transmitted in equal proportions from both the chiral components. To illustrate **10CSA** interactions with $\alpha_2\beta_2$-**P**, a conformational search was performed using the $\alpha_2\beta_2$-**P**[SO_4^{2-}][HSO_4^-]$_4$ structure for building starting geometries. Corresponding noncovalent interactions of the major conformer are illustrated in Figure S14.

In order to calculate e.e. of unknown analyte chiral composition, we have formulated a simple protocol to follow. For example, 19 eq. of an unknown enantiomeric mixture of **10CSA** was complexed with $\alpha_2\beta_2$-**P** (Figure S11). The inner core system shows clearly split resonance signals with $\Delta\sigma$ = 0.13 ppm. A small addition of **10CSA(R)** decreased $\Delta\sigma$ (0.051 ppm), while a similar amount of **10CSA(S)** increased $\Delta\sigma$ (0.224 ppm) revealing the predominant enantiomeric identity (S > R). Next, a linear calibration plot with 7 data points (e.e.$_{10CSA(S)}$ = 0, 48, 58, 69, 79, 90, 100 %) was constructed. It is worth noting that it is possible to generate a calibration curve from a single measured enantiopure point due to the linear dependency with the second point being $\Delta\sigma$ = 0 ppm. where e.e. = 0%. Lastly, the $\Delta\sigma$ can be fitted to the calibration plot and reveal the unknown e.e. (21%) based on the inner N–H chemical shift difference between the split peaks (for more information on this example see figure S12). The same principle can be applied to quickly test the purity of enantiomers, i.e. by equally pre-mixing stereoisomers of interest with opposite chirality, in the event of matching purity, the $\alpha_2\beta_2$-**P** signals will remain isochronous. Overall, monitoring changes of this inner core system in a model chiral environment demonstrates a powerful tool for easy and sensitive detection of enantiomeric compositions.

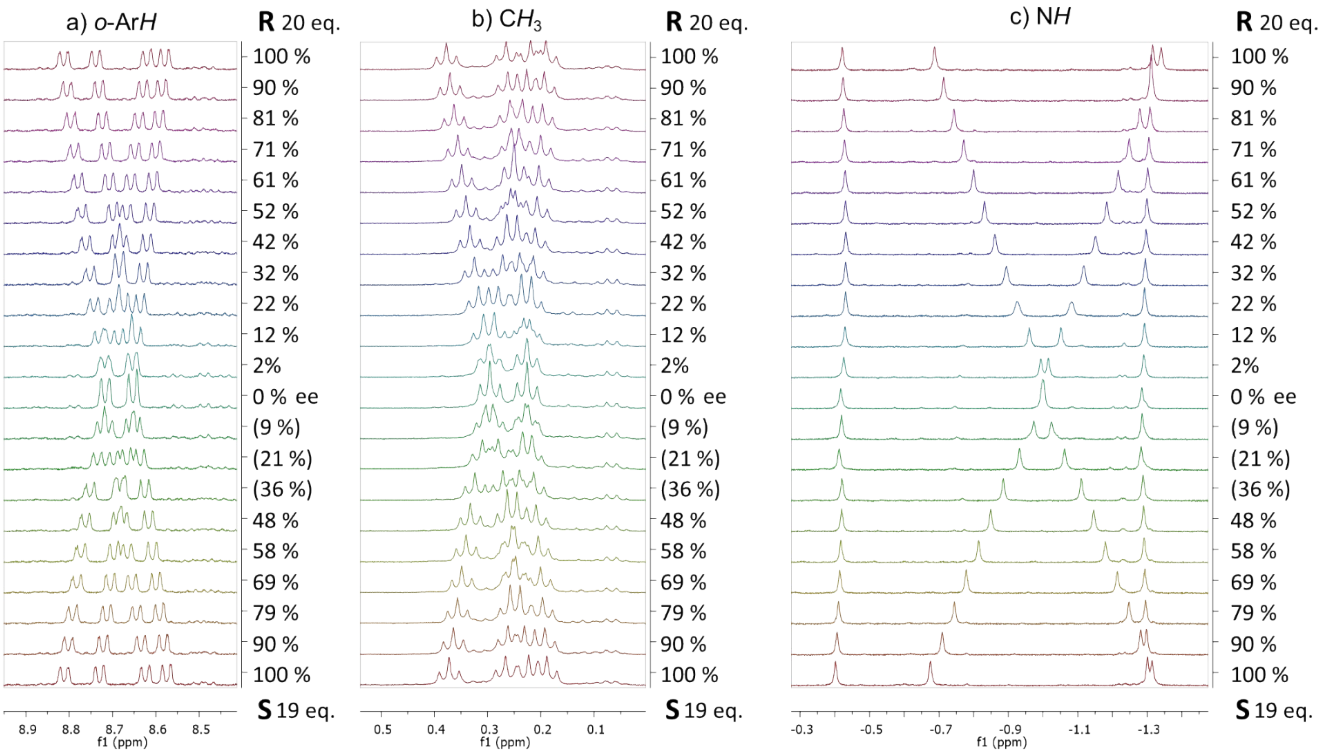


Figure S10. ^{1}H NMR representation of the three non-overlapping regions for the detection of e.e. (values in brackets calculated from the plot see Figure S8-S9: a) *o*-Ar*H*, g) C*H*$_3$, and h) inner core system N–*H*. Two different concentrations of analytes (20 eq. of **10CSA(R)** and 19 eq. of **10CSA(S)**) were used to highlight the diversity of the enantiomeric excess detection using $\alpha_2\beta_2$-**P**. Spectra recorded in CD_3CN.

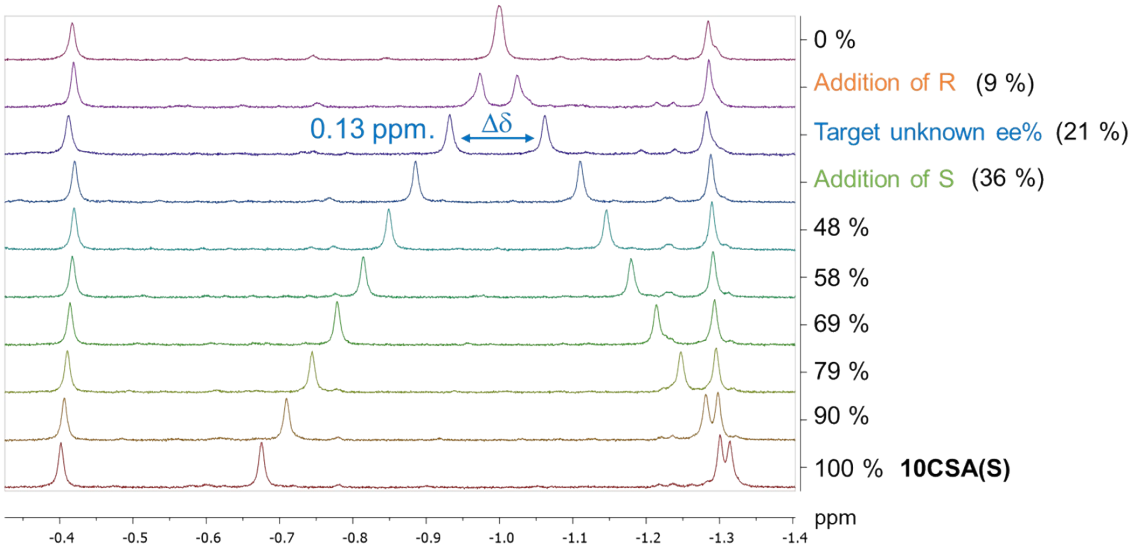
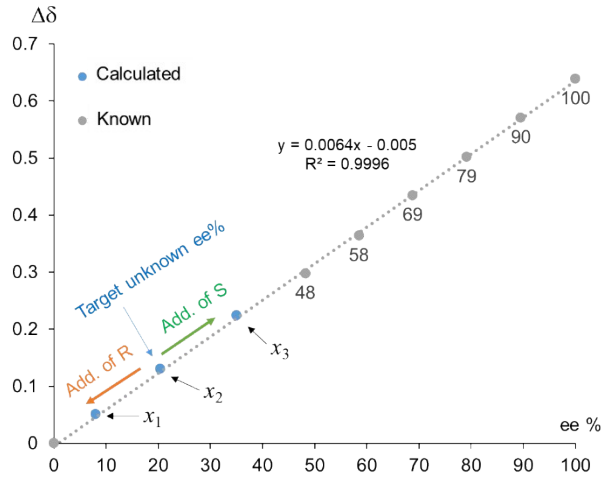


Figure S11. The ^{1}H NMR spectra of the inner core system N–*H* unknown e.e.% target compound (highlighted in blue), after small **10CSA(R)** (red) and **10CSA(S)** (green) additions, and the rest of the spectra for the construction of the calibration curve. Spectra recorded in CD_3CN.



From the calibrated curve:

$$y = 0.0064 \cdot x - 0.005$$

$$x(ee) = \frac{y(\Delta\delta) + 0.005}{0.0064}$$

$$x_1 = \frac{0.051 + 0.005}{0.0064} = 8.75 \approx 9\%$$

$$x_2 = \frac{0.130 + 0.005}{0.0064} = 21.09 \approx 21\%$$

$$x_3 = \frac{0.224 + 0.005}{0.0064} = 35.78 \approx 36\%$$

One point system:

$$x(ee) = \frac{y(\Delta\delta) \cdot 100}{y(\Delta\delta_{max})}$$

$$x_1 = \frac{0.051 \cdot 100}{0.639} = 7.98 \approx 8\%$$

$$x_2 = \frac{0.130 \cdot 100}{0.639} = 20.34 \approx 20\%$$

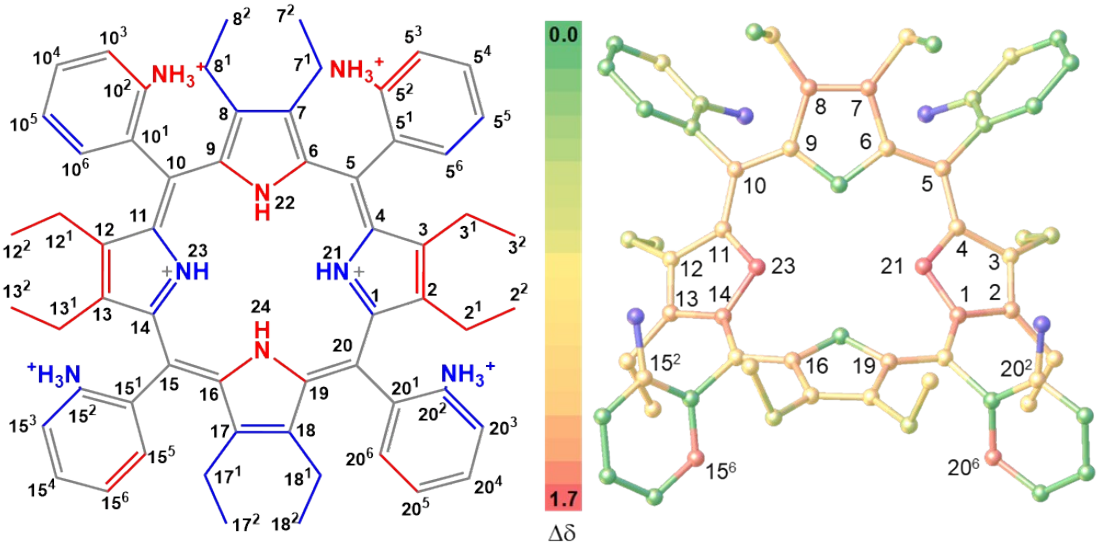
$$x_3 = \frac{0.224 \cdot 100}{0.639} = 35.05 \approx 35\%$$

Figure S12. Graph of the Δσ dependence on the ee% with 19eq. of **10CSA**. Measured from the ^{1}H NMR split N–*H* signals recorded in CD_3CN (at 100 ee% — 19 eq. $\alpha_2\beta_2$-**P·10CSA(S)**). On the right side: calculations of the unknown ee from the calibrated curve and using a one-point system. Note, calculations using a one-point system should only be used for quick, approximate determinations of the ee.

^{13}C and ^{15}N Investigations

To further understand the transfer of chirality to the atropisomeric receptor systems **P**, we have performed 2D NMR analyses with enantiopure **10CSA**. The ^{15}N resonance signals obtained from ^{1}H–^{15}N HSQC varied from 125 to 135 ppm and correlated well with the corresponding inner core system protons. The $\Delta\sigma_{max}$ of the ^{15}N were found to be $\alpha_2\beta_2$-**P** (1.67 ppm) > $\alpha\beta\alpha\beta$-**P** (1.03 ppm) > $\alpha_3\beta$-**P** (0.77 ppm) > α_4-**P** (0 ppm) (Table S2). Full 2D NMR analyses to determine all of the resonance signals were done for $\alpha_2\beta_2$-**P**, $\alpha\beta\alpha\beta$-**P** and α_4-**P** with **10CSA**(**S**). Unfortunately, due to formation of diastereomers in $\alpha_3\beta$-**P·10CSA**(**S**) the large number of signals hampered detailed NMR analysis (Figure S59-S64). While some of the ^{1}H resonance signals in α_4-**P·10CSA**(**S**) showed possible splitting patterns, the $\Delta\sigma_{max}$ were found to not surpass 0.08 ppm. In $\alpha_2\beta_2$-**P·10CSA**(**S**) and $\alpha\beta\alpha\beta$-**P·10CSA**(**S**) other than inner core N–*H* the ^{1}H $\Delta\sigma_{max}$ showed to be highest in *o*-Ar-*H* (0.16 – 0.2 ppm) and certain C*H*$_3$ groups (0.22 ppm in $\alpha\beta\alpha\beta$-**P** 2^2 and 12^2, 7^2 and 17^2; 0.16 ppm in $\alpha_2\beta_2$-**P** 2^2 and 13^2 positions). Only particular ^{13}C in the $\alpha\beta\alpha\beta$-**P·10CSA**(**S**) 24-atom macrocycle displayed $\Delta\sigma_{max}$ with >0.11 ppm (0.47 ppm between 6 and 16, 1 and 11; 0.8 ppm between 8 and 18, 3 and 13 positions). Similarly, in α_4-**P·10CSA**(**S**) only between 1 and 4, 13 and 12 positions the ^{13}C $\Delta\sigma_{max}$ (0.23 ppm) was observed to be >0.11 ppm. On the other hand, the majority of the $\alpha_2\beta_2$-**P·10CSA**(**S**) pyrrolic ^{13}C displayed $\Delta\sigma_{max}$ >0.2 ppm, reaching 0.98 ppm between 1 and 14 positions. In the phenyl rings, except for the previously discussed significant ^{13}C $\Delta\sigma_{max}$ (1.3 ppm) between 15^6 and 20^6 in $\alpha_2\beta_2$-**P·10CSA**(**S**), other phenyl positions and likewise in other atropisomeric species displayed $\Delta\sigma_{max}$ <0.31 ppm.

Table S1. Comparison of ^{13}C and ^{15}N atom resonances $\alpha_2\beta_2$-**P·BSA**,[20] $\alpha_2\beta_2$-**P·10CSA**(**S**) and $\alpha_2\beta_2$-**P·10CSA**(**SR**) complexes. The $\Delta\sigma_{max}$ (ppm) in ^{13}C and ^{15}N atoms of $\alpha_2\beta_2$-**P·10CSA**(**S**) recorded with 20 eq. of analyte. The signal positions were determined using various NMR techniques (CD_3CN). Top left side, illustration of the $\alpha_2\beta_2$-**P** (blue — above and red — below the plane) with corresponding positions. The highlighted positions in the illustration on the top right side shows $\Delta\sigma \geq 0.3$ ppm (atoms in blue are peripheral nitrogen atoms that did not resonate).



Pos.	BSA	CSA-SR	CSA-S
		δ C or δ N	
1	144.1	144.1	143.7
2	142.0	142.1	142.3
2^1	19.1	19.3	19.5
2^2	15.9	16.0	16.0
3	143.9	143.4	143.2
3^1	20.5	20.6	20.5
3^2	15.2	15.6	15.5
4	143.9	144.3	144.5
5	114.9	114.1	113.9
5^1	132.1	132.0	132.0
5^2	133.1	133.1	133.1
5^3	130.6	130.5	130.6
5^4	134.2	133.8	133.8
5^5	126.4	125.9	125.9
5^6	139.8	141.0	141.1
6	142.8	143.0	142.8
7	142.9	141.9	141.4
7^1	19.1	19.4	19.2
7^2	15.9	16.4	16.3
8	142.9	141.9	142.3
8^1	19.1	19.4	19.4
8^2	15.9	16.4	16.3
9	142.8	143.0	143.1
10	114.9	114.1	114.3
10^1	132.1	132.0	131.9
10^2	133.1	133.1	133.2
10^3	130.6	130.5	130.4
10^4	134.2	133.8	133.8
10^5	126.4	125.9	125.9
10^6	139.8	141.0	141.1
11	143.9	144.3	144.1
12	143.9	143.4	143.4

Pos.	BSA	CSA-SR	CSA-S
		δ C or δ N	
12^1	20.5	20.6	20.6
12^2	15.2	15.6	15.6
13	142.0	142.1	141.7
13^1	19.1	19.3	19.3
13^2	15.9	16.0	15.8
14	144.1	144.1	144.7
15	112.0	112.6	112.5
15^1	132.0	132.3	132.3
15^2	133.0	134.0	134.3
15^3	130.5	129.9	129.7
15^4	134.2	133.8	133.8
15^5	125.9	126.3	126.1
15^6	140.4	139.3	139.9
16	144.9	145.6	145.9
17	145.6	144.4	144.4
17^1	20.8	20.6	20.6
17^2	15.4	15.4	15.3
18	145.6	144.4	144.2
18^1	20.8	20.6	20.5
18^2	15.4	15.4	15.5
19	144.9	145.6	145.3
20	112.0	112.6	112.7
20^1	132.0	132.3	132.3
20^2	133.0	134.0	134.6
20^3	130.5	129.9	129.7
20^4	134.2	133.8	133.8
20^5	125.9	126.3	126.1
20^6	140.4	139.3	138.6
21	127.0	127.6	128.5
22	126.9	128.0	128.0
23	127.0	127.6	126.9
24	125.9	131.9	132.2

δ C or δ N	
Pos.	Δδ
1 and 14	0.98
2 and 13	0.59
2^1 and 13^1	0.26
2^2 and 13^2	0.22
3 and 12	0.22
3^1 and 12^1	0.12
3^2 and 12^2	0.12
4 and 11	0.4
6 and 9	0.35
7 and 8	0.88
7^1 and 8^1	0.19
7^2 and 8^2	0.02
10 and 5	0.47
10^1 and 5^1	0.06
10^2 and 5^2	0.09
10^3 and 5^3	0.14
10^4 and 5^4	0
10^5 and 5^5	0
10^6 and 5^6	0.04
15 and 20	0.19
15^1 and 20^1	0
15^2 and 20^2	0.3
15^3 and 20^3	0
15^4 and 20^4	0
15^5 and 20^5	0
15^6 and 20^6	1.3
16 and 19	0.52
17 and 18	0.2
17^1 and 18^1	0.12
17^2 and 18^2	0.18
21 and 23	1.67

Table S2. Comparison of ^{13}C and ^{15}N atom resonances a) α_4-**P·BSA**,[20] α_4-**P·10CSA**(**S**), α_4-**P·10CSA**(**SR**) and, b) $\alpha\beta\alpha\beta$-**P·BSA**,[20] $\alpha\beta\alpha\beta$-**P·10CSA**(**S**), $\alpha\beta\alpha\beta$-**P·10CSA**(**SR**) complexes. The signal positions were determined using various NMR techniques (CD_3CN) recorded with 20 eq. of analyte.

a) α_4-**P**

Pos.	BSA	CSA-SR δ C or δ N	CSA-S	Pos.	BSA	CSA-SR δ C or δ N	CSA-S
1	143.2	143.9	141.9	12^1	18.2	19.2	19.2
2	142.2	145.0	143.7	12^2	16.2	16.1	16.2
2^1	18.2	19.2	19.2	13	142.2	145.0	143.7
2^2	16.2	16.1	16.2	13^1	18.2	19.2	19.2
3	142.2	145.0	143.7	13^2	16.2	16.1	16.1
3^1	18.2	19.2	19.2	14	143.2	143.9	141.8
3^2	16.2	16.1	16.1	15	114.1	113.8	113.8
4	143.2	143.9	141.8	15^1	131.7	131.6	131.6
5	114.1	113.8	113.8	15^2	133.0	133.5	133.4
5^1	131.7	131.6	131.6	15^3	126.2	126.3	126.2
5^2	133.0	133.5	133.4	15^4	134.3	134.1	134.0
5^3	126.2	126.3	126.2	15^5	130.7	130.5	130.5
5^4	134.3	134.1	134.0	15^6	140.6	141.0	140.9
5^5	130.7	130.5	130.5	16	144.9	143.8	144.9
5^6	140.6	141.0	140.9	17	144.8	141.9	143.9
6	144.9	143.8	144.9	17^1	20.4	20.4	19.2
7	144.8	141.9	143.9	17^2	15.0	15.7	15.7
7^1	20.4	20.4	19.2	18	144.8	141.9	143.9
7^2	15.0	15.7	15.7	18^1	20.4	20.4	19.2
8	144.8	141.9	143.9	18^2	15.0	15.7	15.6
8^1	20.4	20.4	19.2	19	144.9	143.8	145.2
8^2	15.0	15.7	15.6	20	114.1	113.8	113.8
9	144.9	143.8	145.2	20^1	131.7	131.6	131.7
10	114.1	113.8	113.8	20^2	133.0	133.5	133.5
10^1	131.7	131.6	131.7	20^3	126.2	126.3	126.3
10^2	133.0	133.5	133.5	20^4	134.3	134.1	134.1
10^3	126.2	126.3	126.3	20^5	130.7	130.5	130.5
10^4	134.3	134.1	134.1	20^6	140.6	141.0	141.1
10^5	130.7	130.5	130.5	21	125.8	127.9	127.7
10^6	140.6	141.0	141.1	22	125.1	125.8	126.1
11	143.2	143.9	141.9	23	125.8	127.9	127.7
12	142.2	145.0	143.7	24	125.1	125.8	126.1

b) $\alpha\beta\alpha\beta$-**P**

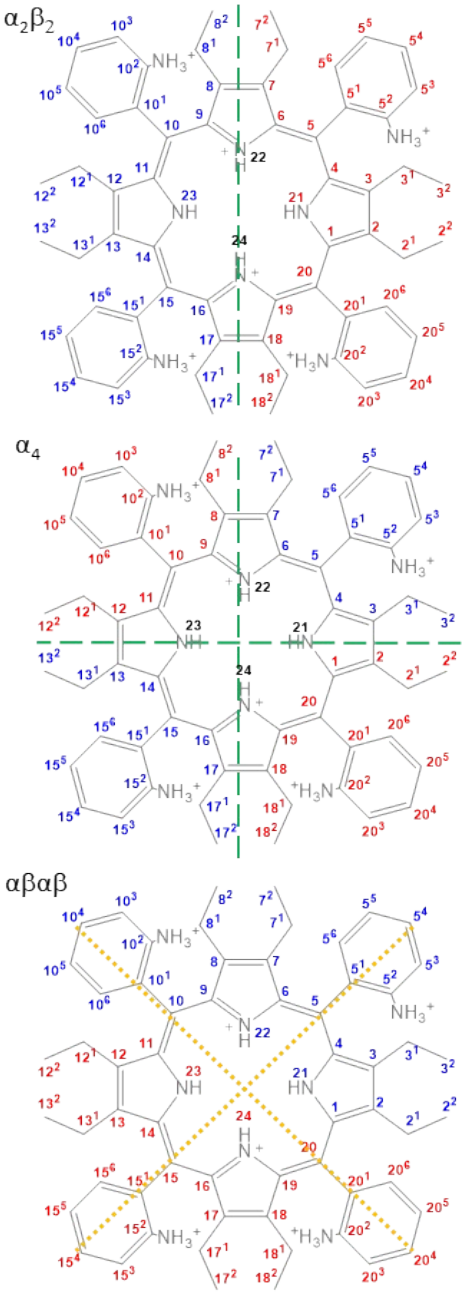
Pos.	BSA	CSA-SR δ C or δ N	CSA-S	Pos.	BSA	CSA-SR δ C or δ N	CSA-S
1	143.6	145.0	145.3	12^1	19.8	20.5	20.2
2	143.8	142.7	142.7	12^2	15.3	15.8	15.8
2^1	19.8	20.5	20.2	13	144.0	142.5	142.2
2^2	15.3	15.8	15.8	13^1	20.6	19.7	19.7
3	144.0	142.5	143.0	13^2	15.6	16.1	16.1
3^1	20.6	19.7	19.7	14	145.0	143.9	143.9
3^2	15.6	16.1	16.1	15	113.1	113.2	113.2
4	145.0	143.9	143.9	15^1	132.1	132.2	132.5
5	113.1	113.2	113.1	15^2	132.9	134.6	133.9
5^1	132.1	132.2	132.2	15^3	126.0	125.8	126.1
5^2	132.9	134.6	133.9	15^4	134.4	133.8	133.9
5^3	126.0	125.8	125.9	15^5	130.9	129.9	130.3
5^4	134.4	133.8	133.9	15^6	140.1	140.3	133.9
5^5	130.9	129.9	130.1	16	143.6	145.0	145.3
5^6	140.1	140.3	133.9	17	143.8	142.7	142.7
6	143.6	145.0	144.8	17^1	19.8	20.5	20.2
7	143.8	142.7	142.8	17^2	15.3	15.8	15.8
7^1	19.8	20.5	20.2	18	144.0	142.5	143.0
7^2	15.3	15.8	15.8	18^1	20.6	19.7	19.7
8	144.0	142.5	142.2	18^2	15.6	16.1	16.1
8^1	20.6	19.7	19.7	19	145.0	143.9	143.9
8^2	15.6	16.1	16.1	20	113.1	113.2	113.1
9	145.0	143.9	143.9	20^1	132.1	132.2	132.2
10	113.1	113.2	113.2	20^2	132.9	134.6	133.9
10^1	132.1	132.2	132.5	20^3	126.0	125.8	125.9
10^2	132.9	134.6	133.9	20^4	134.4	133.8	133.9
10^3	126.0	125.8	126.1	20^5	130.9	129.9	130.1
10^4	134.4	133.8	133.9	20^6	140.1	140.3	133.9
10^5	130.9	129.9	130.3	21	126.5	129.0	129.2
10^6	140.1	140.3	133.9	22	126.5	129.0	128.2
11	143.6	145.0	144.8	23	126.5	129.0	128.2
12	143.8	142.7	142.8	24	126.5	129.0	129.2

Table S3. Comparison of the $\Delta\sigma_{max}$ (ppm) in ^{13}C, ^{15}N and ^{1}H atoms of $\alpha_2\beta_2$-**P·10CSA(S)**, α_4-**P·10CSA(S)**, and $\alpha\beta\alpha\beta$-**P·10CSA(S)** recorded with 20 eq. of analyte. The signal positions were determined using 2D NMR techniques (CD_3CN). On the right side, illustration of the $\alpha_2\beta_2$-**P**, α_4-**P**, and $\alpha\beta\alpha\beta$-**P** with highlighted positions (blue and red represents the splitting signals between them), dashed green line — mirror planes, yellow dotted lines — inversion points.

$\alpha_2\beta_2$-P-10CSA(S)		
Pos.	Δδ C or Δδ N	Δδ H
1 and 14	0.98	
2 and 13	0.59	
2^1 and 13^1	0.26	0.03
2^2 and 13^2	0.22	0.16
3 and 12	0.22	
3^1 and 12^1	0.12	0.09
3^2 and 12^2	0.12	0.03
4 and 11	0.4	
6 and 9	0.35	
7 and 8	0.88	
7^1 and 8^1	0.19	0.03
7^2 and 8^2	0.02	0.07
10 and 5	0.47	
10^1 and 5^1	0.06	
10^2 and 5^2	0.09	
10^3 and 5^3	0.14	
10^4 and 5^4	0	
10^5 and 5^5	0	
10^6 and 5^6	0.04	0.2
15 and 20	0.19	
15^1 and 20^1	0	
15^2 and 20^2	0.3	
15^3 and 20^3	0	
15^4 and 20^4	0	
15^5 and 20^5	0	
15^6 and 20^6	1.3	0.17
16 and 19	0.52	
17 and 18	0.2	
17^1 and 18^1	0.12	0.09
17^2 and 18^2	0.18	0.03
21 and 23	1.67	0.66
22 and 24	0	0

α_4-P-10CSA(S)		
Pos.	Δδ C or Δδ N	Δδ H
1 and 4	0.23	
2 and 3	0.01	
2^1 and 3^1	0	0
2^2 and 3^2	0.11	0.08
6 and 9	0.1	
7 and 8	0.08	
7^1 and 8^1	0	0
7^2 and 8^2	0.03	0
10 and 5	0.06	
10^1 and 5^1	0.05	
10^2 and 5^2	0.18	
10^3 and 5^3	0.07	0.07
10^4 and 5^4	0.06	0.01
10^5 and 5^5	0	0
10^6 and 5^6	0.19	0.01
13 and 12	0.23	
13^1 and 12^1	0.01	
13^2 and 12^2	0	0
11 and 14	0.11	0.08
16 and 19	0.1	
17 and 18	0.08	
17^1 and 18^1	0	0
17^2 and 18^2	0.03	0
15 and 20	0.06	
15^1 and 20^1	0.05	
15^2 and 20^2	0.18	
15^3 and 20^3	0.07	0.07
15^4 and 20^4	0.06	0.01
15^5 and 20^5	0	0
15^6 and 20^6	0.19	0.01
21 and 23	0	0
22 and 24	0	0

$\alpha\beta\alpha\beta$-P-10CSA(S)		
Pos.	Δδ C or Δδ N	Δδ H
1 and 11	0.47	
2 and 12	0.09	
2^1 and 12^1	0	0.02
2^2 and 12^2	0	0.22
4 and 14	0	
3 and 13	0.8	
3^1 and 13^1	0	0.1
3^2 and 13^2	0.03	0.03
6 and 16	0.47	
7 and 17	0.09	
7^1 and 17^1	0	0.02
7^2 and 17^2	0	0.22
9 and 19	0	
8 and 18	0.8	
8^1 and 18^1	0	0.1
8^2 and 18^2	0.03	0.03
15 and 5	0.1	
15^1 and 5^1	0.31	
15^2 and 5^2	0	
15^3 and 5^3	0.16	0.05
15^4 and 5^4	0	0
15^5 and 5^5	0.22	0
15^6 and 5^6	0	0.16
10 and 20	0.1	
10^1 and 20^1	0.31	
10^2 and 20^2	0	
10^3 and 20^3	0.16	0.05
10^4 and 20^4	0	0
10^5 and 20^5	0.22	0
10^6 and 20^6	0	0.16
21 and 23	1.03	0.52
22 and 24	1.03	0.52



SUPPORTING INFORMATION

Structural Determination of $\alpha_2\beta_2$-**P**$[SO_4^{2-}][HSO_4^-]_4$

Table S4: Details of XRD data refinement of $\alpha_2\beta_2$-**P**$[SO_4^{2-}][HSO_4^-]_4$

Compound	$\alpha_2\beta_2$-**P**$[SO_4^{2-}][HSO_4^-]_4$
Internal code	KN007
CCDC #	2143572
Empirical formula	$C_{60}H_{84.38}N_8O_{23.95}S_5$
Formula weight	1461.23
Temperature/K	100(2)
Crystal system	Monoclinic
Space group	$P2_1/n$
a/Å	18.7594(9)
b/Å	16.0888(8)
c/Å	24.9988(12)
α/°	90
β/°	90.962(2)
γ/°	90
Volume/Å³	7544.0(6)
Z	4
D_{calc} *g/cm³*	1.287
μ/mm⁻¹	2.066
F(000)	3088.0
Crystal size/mm³	0.15 × 0.1 × 0.07
Radiation	CuKα
Wavelength/Å	1.54178
2θ/°	5.844 to 140.326
Reflections collected	94031
Independent reflections	14243
R_{int}	0.0578
R_{sigma}	0.0345
Restraints	3307
Parameters	1354
GooF	1.104
R_1 *[I> 2σ (I)]*	0.1354
wR_2 *[I> 2σ (I)]*	0.3912
R_1 *[all data]*	0.1537
wR_2 *[all data]*	0.4253
Largest peak/e Å⁻³	1.73
Deepest hole/e Å⁻³	-0.99

basis	Δ_{ip}	δ_{ip}	B_{2g}	B_{1g}	$E_u(x)$	$E_u(y)$	A_{1g}	A_{2g}
min.	0.75	0.18	-0.02	0.11	-0.02	0.00	-0.74	0.01
ext.	0.88	0.13	-0.02	0.11	-0.02	0.00	-0.74	0.01
			0.01	0.01	-0.03	0.01	-0.46	0.01
total	1.11	0.00	-0.02	0.11	-0.02	0.00	-0.71	0.01
			0.01	0.01	-0.03	0.01	-0.45	0.01
			0.01	-0.01	-0.01	-0.01	0.71	0.00
			0.00	0.00	-0.02	0.02	0.00	0.00
			0.01	0.01	0.01	-0.01	-0.07	0.00
			0.00	0.00	0.01	-0.01	-0.07	
					-0.01	0.02		
					0.00	0.00		
					0.00	-0.02		
					0.00	0.00		
					0.00	0.01		
comp.	1.11	0.00	0.02	0.11	0.04	0.04	1.10	0.02

basis	Δ_{oop}	δ_{oop}	B_{2u}	B_{1u}	A_{2u}	$E_g(x)$	$E_g(y)$	A_{1u}
min.	3.88	0.11	-3.88	-0.02	0.02	0.02	-0.01	0.02
ext.	3.92	0.00	-3.87	-0.02	0.02	0.02	-0.01	0.02
			0.65	-0.01	-0.02	-0.04	0.00	0.00
total	3.92	0.00	-3.87	-0.02	0.02	0.02	-0.01	0.02
			0.65	-0.01	-0.02	-0.04	0.00	0.00
			0.07	-0.01	0.00	0.01	-0.03	
						-0.01	0.00	
						-0.01	0.00	
comp.	3.92	0.00	3.92	0.02	0.03	0.04	0.03	0.02

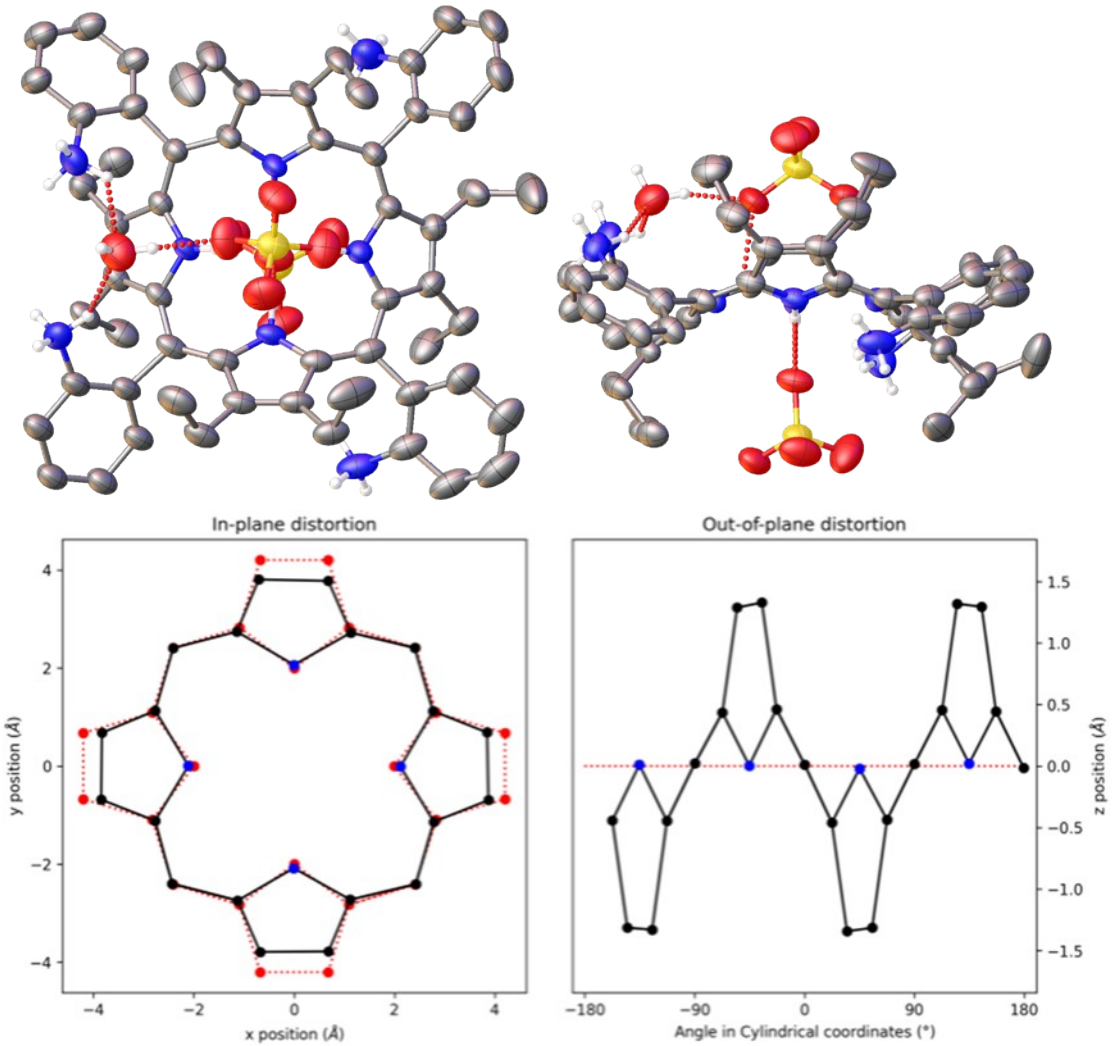


Figure S13. Left: in-plane (ip) and out-of-plane (oop) NSD results of $\alpha_2\beta_2$-**P**[SO_4^{2-}][HSO_4^-]$_4$; Right bottom: out-of-plane and in-plane skeletal plots of the porphyrin $\alpha_2\beta_2$-**P**[SO_4^{2-}][HSO_4^-]$_4$ core. Porphyrin $\alpha_2\beta_2$-**P** is represented in black(C) and blue(N), with the reference structure (CuTPP) in red dotted lines.[2] Right top: top view and sideview of $\alpha_2\beta_2$-**P**-SO_4 structure. Non-essential hydrogens, majority of counter anions and solvent molecules omitted for clarity, thermal ellipsoids shown at 50% probability.

Computational Analysis

Geometry analysis

Based on the crystallographic data we concluded that one porphyrin cation binds two camphorsulfonic acid molecules. Using this ratio and the $\alpha_2\beta_2$-**P**[SO_4^{2-}][HSO_4^-]$_4$ crystal data the corresponding host-guest complexes were built with subsequent conformation search. According to the Boltzmann distribution, in acetonitrile one major conformer **A** (90%) and one minor conformer **B** (up to 10%) being higher in energy by 1.30 kcal/mol are presented (Table S5).

In all host-guest complexes, the porphyrin plane is significantly distorted (the C_β-C_β-$C_{\beta opp}$-$C_{\beta opp}$ angles varying in the range of 0˚- 25˚ and N-C_α-C_α-N angles varying in the range of 37˚- 48˚) because of a steric hindrance between the peripheral substituents (for NSD profile check Figure S13). This deformation results in appearance of two cavities on both sides of the porphyrin macrocycle, which differ by the position of NH_3^+ groups. In one cavity the ammonia groups are placed on the same side of the cavity and in another – on the opposite sides (Figure S14a), and the distance between two nitrogen atoms increases from 4.899 Å to 8.7174 Å, respectively. The difference in the position of NH_3^+ groups results in a non-identical mode of the binding of two camphorsulfonic acids (“standing” and “lying”). Both the NBO and AIMALL analysis showed that the guest molecules interact with the porphyrin cation through the formation of four H-bonds, which cause significant elongation of the N–H bonds by 0.08-0.03 Å in the case of NH_3^+ groups and by 0.03 Å in the case of inner core protons for the major conformer **A** (Figure S14b, Tables S6 and S7).

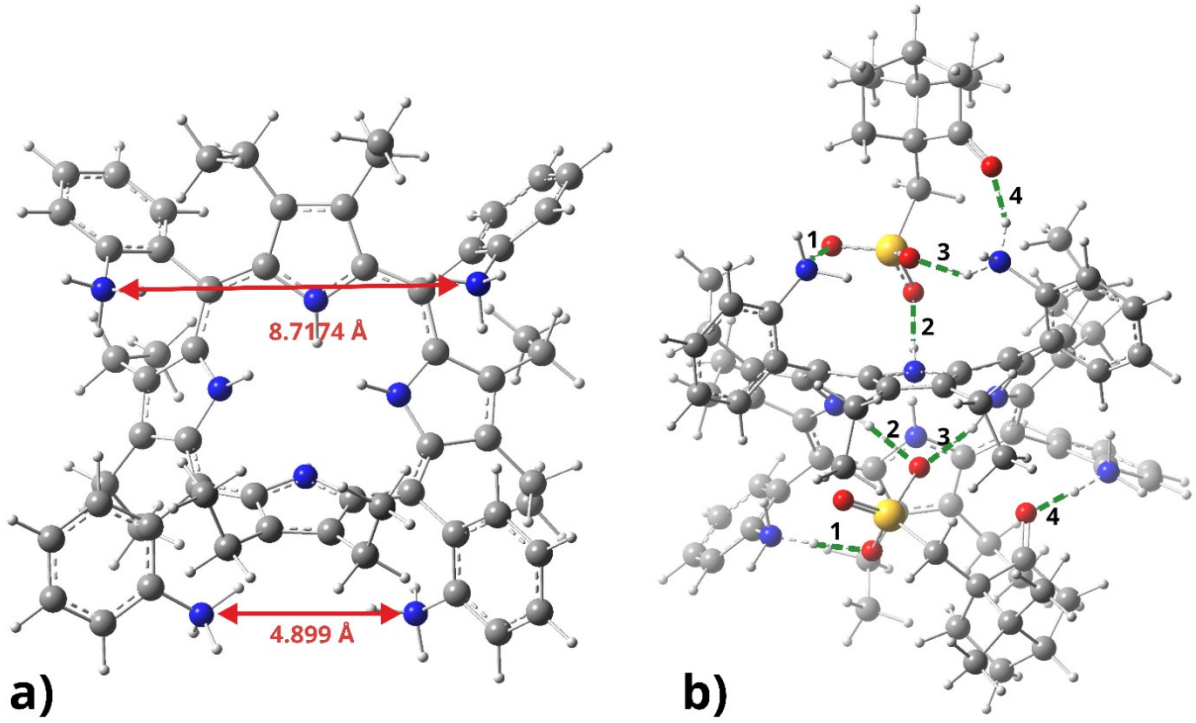


Figure S14. a) Position of NH_3^+ groups in porphyrin cation; **b)** major conformer **A** of the $\alpha_2\beta_2$-**P·10CSA(R)** complex**,** where hydrogen bonds formed between porphyrin and two guest molecules are shown by the green dashed lines and numbered.

Table S5. Relative energies and Boltzmann distribution of $\alpha_2\beta_2$-**P·10CSA(R)** conformers.

Conf.	El. energy, Hartree (def2-SVP)	El. energy, Hartree (def2-SVP)	Gibbs Free energy correction, Hartree	Relative Gibbs Free energy, kcal/mol	Boltzmann distribution, %
A	-4942,80946	-4947,6688	1,534213	0	90
B	-4942,80783	-4947,66664	1,534117	1,3	10
C	-4942,79772	-4947,65773	1,533141	6,28	0
D	-4942,79733	-4947,6544	1,534908	9,48	0
E	-4942,79347	-4947,65221	1,532854	9,56	0
F	-4942,79077	-4947,64779	1,536709	14,75	0
G	-4942,77487	-4947,63394	1,532741	20,96	0
H	-4942,7644	-4947,6243	1,532105	26,6	0

Table S6. Perturbation theory energy analysis of A conformer of $\alpha_2\beta_2$-**P·10CSA(R)** complex.

Num. of H-bond	"standing" guest			"lying" guest		
	Donor NBO	**Acceptor NBO**	**E(2), kcal/mol**	**Donor NBO (I)**	**Acceptor NBO (j)**	**E(2), kcal/mol**
1	BD (1) O 138 - S 165	RY*(2) H 103	0.12	BD (1) N 127 - H 130	RY*(3) O81	0.06
	BD (1) O 138 - S 165	RY*(3) H 103	0.1	BD (1) N 127 - H 130	RY*(9) O81	0.06
	BD (1) O 138 - S 165	RY*(4) H 103	0.1	BD (1) N 127 - H 130	BD*(1) O81 - S 169	0.52
	BD (1) O 138 - S 165	RY*(5) H 103	0.07	BD (1) N 127 - H 130	RY*(3) O81	1.42
	BD (1) O 138 - S 165	BD*(1) N 100 - H 103	0.31	BD (1) O81 - S 169	RY*(2) H 130	0.09
	CR (1) O 138	BD*(1) N 100 - H 103	0.84	BD (1) O81 - S 169	RY*(3) H 130	0.09
	LP (1) O 138	RY*(1) H 103	0.29	BD (1) O81 - S 169	RY*(4) H 130	0.18
	LP (1) O 138	RY*(3) H 103	0.14	BD (1) O81 - S 169	BD*(1) N 127 - H 130	0.48
	LP (1) O 138	BD*(1) N 100 - H 103	10.53	CR (1) O81	RY*(1) H 130	0.06
	LP (2) O 138	RY*(6) H 103	0.15	CR (1) O81	RY*(2) H 130	0.06
	LP (3) O 138	RY*(1) H 103	1.08	CR (1) O81	BD*(1) N 127 - H 130	1.24
	LP (3) O 138	RY*(2) H 103	1.28	CR (1) O81	RY*(1) H 130	0.34
	LP (3) O 138	RY*(3) H 103	0.31	CR (1) O81	RY*(2) H 130	0.06
	LP (3) O 138	RY*(4) H 103	0.25	CR (1) O81	RY*(3) H 130	0.06
	LP (3) O 138	RY*(5) H 103	0.08	CR (1) O81	BD*(1) N 127 - H 130	11.71
	LP (3) O 138	RY*(8) H 103	0.07	LP (2) O81	RY*(3) H 130	0.1
	LP (3) O 138	BD*(1) N 100 - H 103	33.97	LP (2) O81	RY*(6) H 130	0.07
	BD*(1) O 138 - S 165	BD*(1) N 100 - H 103	4.38	LP (2) O81	RY*(7) H 130	0.05
	BD (1) N 100 - H 103	RY*(2) O 138	0.07	LP (2) O81	BD*(1) N 127 - H 130	0.32
	BD (1) N 100 - H 103	BD*(1) O 138 - S 165	0.4	LP (3) O81	RY*(1) H 130	0.65
				LP (3) O81	RY*(2) H 130	2.32
				LP (3) O81	RY*(3) H 130	0.13
				LP (3) O81	RY*(4) H 130	0.23
				LP (3) O81	RY*(5) H 130	0.16
				LP (3) O81	RY*(8) H 130	0.06
				LP (3) O81	BD*(1) N 127 - H 130	43.13
				BD*(1) O81 - S 169	BD*(1) N 127 - H 130	5.51
	Total		**54.54**	**Total**		**69.16**
2	BD (1) S 165 - O 167	RY*(2) H53	0.07	BD (1) N50 - H51	RY*(2) O87	0.11

	BD (1) S 165 - O 167	RY*(4) H53	0.05	BD (1) N50 - H51	RY*(3) O87	0.09
	BD (1) S 165 - O 167	BD*(1) N52 - H53	0.35	BD (1) N50 - H51	BD*(1) O87 - S 169	0.18
	CR (1) O 167	BD*(1) N52 - H53	0.56	BD (1) O87 - S 169	RY*(3) H51	0.16
	LP (1) O 167	RY*(1) H53	0.12	BD (1) O87 - S 169	RY*(6) H51	0.11
	LP (1) O 167	RY*(2) H53	0.33	BD (1) O87 - S 169	BD*(1) N50 - H51	0.69
	LP (1) O 167	RY*(3) H53	0.09	CR (1) O87	BD*(1) N50 - H51	0.71
	LP (1) O 167	BD*(1) N52 - H53	8.72	LP (1) O87	RY*(1) H51	0.12
	LP (2) O 167	RY*(2) H53	0.13	LP (1) O87	RY*(2) H51	0.26
	LP (2) O 167	BD*(1) N52 - H53	0.88	LP (1) O87	BD*(1) N50 - H51	8.05
	LP (3) O 167	RY*(1) H53	0.05	LP (2) O87	RY*(7) H51	0.07
	LP (3) O 167	RY*(2) H53	2.3	LP (3) O87	RY*(1) H51	2.29
	LP (3) O 167	RY*(3) H53	0.61	LP (3) O87	RY*(2) H51	0.45
	LP (3) O 167	RY*(4) H53	0.09	LP (3) O87	RY*(3) H51	0.08
	LP (3) O 167	RY*(5) H53	0.46	LP (3) O87	RY*(4) H51	0.14
	LP (3) O 167	BD*(1) N52 - H53	16.92	LP (3) O87	RY*(5) H51	0.59
	BD*(1) S 165 - O 167	BD*(1) N52 - H53	1.36	LP (3) O87	BD*(1) N50 - H51	18.95
	BD (1) N52 - H53	RY*(1) O 167	0.15	BD*(1) O87 - S 169	BD*(1) N50 - H51	0.8
	BD (1) N52 - H53	RY*(2) O 167	0.06			
	BD (1) N52 - H53	BD*(1) S 165 - O 167	0.18			
	Total		**33.48**	**Total**		**33.85**
3	BD (1) S 165 - O 166	RY*(1) H91	0.36	BD (1) N54 - H55	RY*(1) O87	0.16
	BD (1) S 165 - O 166	RY*(3) H91	0.24	BD (1) N54 - H55	RY*(4) O87	0.06
	BD (1) S 165 - O 166	RY*(4) H91	0.06	BD (1) N54 - H55	BD*(1) O87 - S 169	1.3
	BD (1) S 165 - O 166	BD*(1) N90 - H91	0.08	BD (1) O87 - S 169	RY*(1) H55	1.57
	CR (1) O 166	BD*(1) N90 - H91	0.48	BD (1) O87 - S 169	RY*(2) H55	0.22
	LP (1) O 166	RY*(1) H91	0.52	BD (1) O87 - S 169	RY*(3) H55	0.63
	LP (1) O 166	RY*(2) H91	0.21	BD (1) O87 - S 169	BD*(1) N54 - H55	0.06
	LP (1) O 166	RY*(5) H91	0.06	CR (1) O87	BD*(1) N54 - H55	0.48
	LP (1) O 166	BD*(1) N90 - H91	9.85	LP (1) O87	RY*(1) H55	1.93
	LP (2) O 166	BD*(1) N90 - H91	0.46	LP (1) O87	RY*(2) H55	0.14
	LP (3) O 166	RY*(1) H91	2.14	LP (1) O87	RY*(3) H55	0.06
	LP (3) O 166	RY*(3) H91	0.21	LP (1) O87	BD*(1) N54 - H55	15.73
	LP (3) O 166	RY*(5) H91	0.12	LP (2) O87	RY*(1) H55	0.08
	LP (3) O 166	BD*(1) N90 - H91	13.33	LP (2) O87	RY*(4) H55	0.08
	BD*(1) S 165 - O 166	BD*(1) N90 - H91	3.44	LP (2) O87	BD*(1) N54 - H55	0.4
	BD (1) N90 - H91	RY*(1) O 166	0.27	LP (3) O87	RY*(1) H55	0.55
	BD (1) N90 - H91	BD*(1) S 165 - O 166	0.84	LP (3) O87	RY*(6) H55	0.08
				LP (3) O87	BD*(1) N54 - H55	1.08
				BD*(1) O87 - S 169	RY*(3) H55	2.82
				BD*(1) O87 - S 169	BD*(1) N54 - H55	5.06
	Total		**32.07**	**Total**		**32.49**
4	BD (1) C1 - O25	RY*(1) H 147	0.49	BD (1) N66 - H67	BD*(1) C 170 - O 195	0.63
	BD (1) C1 - O25	RY*(2) H 147	0.14	BD*(1) N66 - H67	BD*(1) C 170 - O 195	1.11
	BD (1) C1 - O25	RY*(3) H 147	0.21	BD (1) C 170 - O 195	RY*(5) H67	0.07
	BD (1) C1 - O25	BD*(1) N90 - H 147	0.53	BD (1) C 170 - O 195	BD*(1) N66 - H67	0.59
	BD (2) C1 - O25	RY*(1) H 147	0.1	BD (1) C 170 - O 195	RY*(6) H67	0.08
	BD (2) C1 - O25	RY*(2) H 147	0.16	CR (1) O 195	BD*(1) N66 - H67	0.48

	BD (2) C1 - O25	RY*(3) H 147	0.37	LP (1) O 195	RY*(1) H67	0.21
	BD (2) C1 - O25	RY*(4) H 147	0.07	LP (1) O 195	RY*(4) H67	0.12
	BD (2) C1 - O25	BD*(1) N90 - H 147	0.77	LP (1) O 195	BD*(1) N66 - H67	6.81
	CR (1) O25	BD*(1) N90 - H 147	0.3	LP (2) O 195	RY*(2) H67	1.46
	LP (1) O25	RY*(2) H 147	0.96	LP (2) O 195	RY*(4) H67	0.35
	LP (1) O25	RY*(3) H 147	0.08	LP (2) O 195	RY*(7) H67	0.05
	LP (1) O25	BD*(1) N90 - H 147	17.28	LP (2) O 195	RY*(8) H67	0.06
	LP (2) O25	RY*(2) H 147	0.44	LP (2) O 195	BD*(1) N66 - H67	31.56
	LP (2) O25	BD*(1) N90 - H 147	5.58	BD*(2) C 170 - O 195	BD*(1) N66 - H67	0.09
	BD*(2) C1 - O25	BD*(1) N90 - H 147	0.73			
	BD (1) N90 - H 147	RY*(2) O25	0.23			
	BD (1) N90 - H 147	RY*(4) O25	0.1			
	BD (1) N90 - H 147	BD*(1) C1 - O25	1.01			
	BD (1) N90 - H 147	BD*(2) C1 - O25	0.72			
	Total		**30.27**	**Total**		**43.67**

Table S7. Topological analysis of the electron density of A conformer of $\alpha_2\beta_2$-**P·10CSA(R)** complex.

Guest	Atoms	Electron delocalization index	Electron Dnsity (ρ(r)), a.u.	Laplacian of Rho
"Standing"	O_{138} & H_{103}	3.934	0.06974	0.15993
	O_{167} & H_{53}	4.277	0.04978	0.16467
	O_{166} & H_{91}	3.915	0.04906	0.15993
	O_{25} & H_{147}	4.013	0.05257	0.17475
"Lying"	O_{81} & H_{130}	3.888	0.07823	0.14315
	O_{87} & H_{51}	4.225	0.04127	0.15588
	O_{87} & H_{55}	4.062	0.04127	0.15757
	O_{195} & H_{67}	3.938	0.04127	0.14724

NMR calculations and analysis

In addition, the ^{13}C and ^{15}N NMR shielding tensors and chemical shifts, as well as $\Delta\sigma_{max}$ between the corresponding chemical shifts of the main dominant conformer **A** of $\alpha_2\beta_2$-**P·10CSA(R)** were calculated (Tables S8-S10). The calculated ^{13}C and ^{15}N NMR chemical shifts were found to deviate from the experimentally measured data, while the calculated and experimental $\Delta\sigma_{max}$ values are in good agreement (Table S8), except the $\Delta\sigma_{max}$ between carbons 2 and 13, 3 and 12, and nitrogens 22 and 24, which showed strong deviation (Table S8). The disagreement between these three $\Delta\sigma_{max}$ values can be explained by the presence and influence of solvent molecules situated in a larger cavity. In contrast to the smaller cavity, where only one "lying" guest molecule can be placed, in the larger cavity in addition to the "standing" guest molecule several solvent molecules can be placed as well. However, the continuum solvent model (SND) used represents solvent as a continuous medium and does not count for individual effects of "explicit" molecules, such as H-bonds formation or charge transfer. We suppose that the deviation from the experiment is due to not accounting for the individual solvent molecules' effects in the larger porphyrin cavity. Between other $\Delta\sigma_{max}$ calculated using theoretical and experimental data, a good correlation is observed, proving that the main calculated conformation corresponds to the

dominant conformation presented in solution, which shows non-equivalency of the carbon and nitrogen atoms. To improve agreement with experimental measurements the conformer **A** of $\alpha_2\beta_2$-**P·10CSA(R)** with one and two additional acetonitrile molecules was modelled, and its ^{13}C and ^{15}N NMR shielding tensors and chemical shifts, as well as $\Delta\sigma_{max}$ between the corresponding chemical shifts were calculated. Inclusion of explicit solvent molecules in the model system improved agreement with the experimental data, especially in the case of the $\Delta\sigma_{max}$ between carbons 2 and 13, 3 and 12, and nitrogens 22 and 24 (Table S8). However, the model system is sensitive to the presence of solvent molecules, especially in the case of peripheral atoms. Thus, to perform accurate NMR calculations, the whole solvent shell should be modelled which is beyond the scope of this study.

Table S8. Calculated shielding tensors and chemical shifts of ^{13}C of conformer **A** of $\alpha_2\beta_2$-**P·10CSA(R)** with and without solvent molecules.

Number of atom (in xyz)	$\alpha_2\beta_2$-P·10CSA(R)		$\alpha_2\beta_2$-P·10CSA(R) + 1 acetonitrile		$\alpha_2\beta_2$-P·10CSA(R) + 2 acetonitrile	
	Shielding tensor, ppm	Scaled chemical shift, ppm	Shielding tensor, ppm	Scaled chemical shift, ppm	Shielding tensor, ppm	Scaled chemical shift, ppm
1	-60,94	236,3	-62,83	238,1	-62,84	238,1
2	116,32	63,0	115,74	63,5	115,62	63,7
3	134,16	45,5	133,94	45,7	133,87	45,8
4	132,32	47,3	132,12	47,5	132,25	47,4
7	123,17	56,3	122,69	56,7	123,21	56,2
8	151,34	28,7	151,37	28,7	151,04	29,0
12	161,76	18,5	162,02	18,3	160,60	19,7
16	160,08	20,2	159,39	20,9	160,69	19,6
20	120,29	59,1	119,34	60,0	119,56	59,8
22	151,16	28,9	151,74	28,3	151,49	28,6
26	29,40	147,9	26,32	151,0	26,31	151,0
27	25,99	151,3	27,38	149,9	28,09	149,2
28	32,82	144,6	34,39	143,1	34,25	143,2
29	31,22	146,2	29,68	147,7	27,94	149,4
30	32,35	145,1	31,80	145,6	31,94	145,5
31	58,80	119,2	58,84	119,2	59,68	118,3
32	21,06	156,1	24,13	153,1	25,61	151,6
33	19,13	158,0	24,03	153,2	24,98	152,3
34	31,12	146,3	32,41	145,0	33,27	144,2
35	56,16	121,8	57,43	120,5	57,68	120,3
36	32,74	144,7	31,95	145,5	32,89	144,5
37	25,41	151,9	27,73	149,6	28,54	148,8
38	28,52	148,8	32,52	144,9	33,09	144,3
39	34,09	143,4	32,57	144,8	32,93	144,5
40	27,04	150,3	26,72	150,6	25,79	151,5
41	32,40	145,0	30,66	146,7	27,61	149,7
42	65,21	112,9	63,93	114,2	63,58	114,5
43	63,20	114,9	62,58	115,5	63,69	114,4
46	23,15	154,1	22,41	154,8	22,83	154,4
47	156,80	23,4	156,72	23,5	157,14	23,1
56	159,95	20,3	159,97	20,3	160,12	20,1
60	44,40	133,3	43,94	133,7	43,38	134,3

61	41,49	136,1	41,21	136,4	41,72	135,9
62	49,57	128,2	50,24	127,6	50,81	127,0
63	39,15	138,4	39,29	138,3	39,90	137,7
64	34,31	143,1	34,90	142,6	33,32	144,1
70	158,23	22,0	158,15	22,1	157,80	22,4
74	160,38	19,9	160,20	20,1	159,79	20,5
78	157,83	22,4	157,57	22,6	157,61	22,6
82	156,63	23,6	156,30	23,9	156,46	23,7
86	42,39	135,2	43,85	133,8	43,94	133,7
88	35,22	142,3	32,81	144,6	32,98	144,4
93	158,68	21,6	158,88	21,4	158,80	21,4
95	156,68	23,5	156,86	23,3	156,81	23,4
99	157,59	22,6	157,99	22,2	158,11	22,1
104	154,57	25,6	156,19	24,0	156,17	24,0
106	163,44	16,9	160,73	19,5	160,98	19,3
110	157,01	23,2	157,18	23,0	157,48	22,7
113	157,94	22,3	157,97	22,2	157,83	22,4
117	42,89	134,8	42,18	135,5	42,12	135,5
118	41,72	135,9	41,60	136,0	41,09	136,5
119	49,74	128,1	50,10	127,7	50,04	127,8
121	36,67	140,8	37,11	140,4	37,42	140,1
123	41,87	135,8	42,22	135,4	42,65	135,0
125	38,12	139,4	39,35	138,2	40,00	137,6
131	162,15	18,2	161,69	18,6	161,30	19,0
135	156,29	23,9	156,36	23,8	156,56	23,6
139	32,97	144,5	31,01	146,4	31,17	146,2
140	42,73	134,9	42,79	134,9	43,29	134,4
141	48,71	129,1	48,55	129,2	48,55	129,2
143	36,11	141,4	36,42	141,1	36,65	140,9
145	40,72	136,9	41,06	136,5	41,14	136,5
148	41,20	136,4	41,82	135,8	41,65	136,0
149	43,16	134,5	43,54	134,1	43,62	134,0
150	35,43	142,0	32,88	144,5	32,89	144,5
151	51,34	126,5	50,74	127,1	51,00	126,8
152	41,03	136,6	40,46	137,1	40,50	137,1
154	37,23	140,3	37,08	140,4	37,42	140,1
160	160,91	19,4	160,84	19,4	160,85	19,4
170	-64,04	239,3	-63,16	238,4	-62,54	237,8
171	114,95	64,3	114,51	64,7	114,34	64,9
172	134,30	45,4	132,65	47,0	133,12	46,5
173	133,26	46,4	134,01	45,7	133,87	45,8
176	121,90	57,5	122,19	57,2	122,56	56,9
177	152,41	27,7	152,34	27,7	152,55	27,5
181	162,30	18,0	161,69	18,6	161,27	19,0
185	159,21	21,0	158,81	21,4	158,71	21,5
189	121,29	58,1	120,40	59,0	120,18	59,2
192	146,85	33,1	146,10	33,8	146,05	33,9
196	36,05	141,4	36,61	140,9	36,86	140,7
202			59,37	118,6	60,19	117,8
203			180,29	0,4	179,90	0,8
208					54,36	123,5
209					180,40	0,3

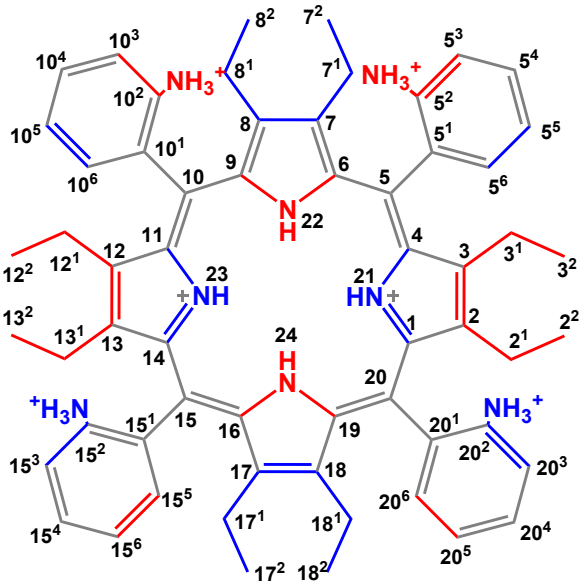
Table S9. Calculated shielding tensors of ^{15}N of conformer **A** of $\alpha_2\beta_2$-**P·10CSA(R)** with and without solvent molecules.

Number of atom (in xyz)	$\alpha_2\beta_2$-P·10CSA(R)		$\alpha_2\beta_2$-P·10CSA(R) + 1 acetonitrile		$\alpha_2\beta_2$-P·10CSA(R) + 2 acetonitrile	
	Shielding tensor, ppm	Scaled chemical shift, ppm	Shielding tensor, ppm	Scaled chemical shift, ppm	Shielding tensor, ppm	Scaled chemical shift, ppm
44	99,30	154,4	87,41	166,3	86,84	166,9
50	99,33	154,4	90,15	163,5	91,46	162,2
52	91,95	161,7	99,64	154,1	100,58	153,1
54	106,38	147,3	101,99	151,7	101,91	151,8
66	180,72	73	180,58	73,1	180,73	73
90	185,44	68,3	184,27	69,4	184,17	69,5
100	173,69	80	173,26	80,4	173,23	80,5
127	178,41	75,3	177,34	76,4	176,61	77,1
			6,94	246,8	2,86	250,8
					-20,51	274,2

Table S10. Calculated and experimentally measured $\Delta\sigma_{max}$ (ppm) between chemical shifts in ^{13}C of of $\alpha_2\beta_2$-**P·10CSA(R)** with and without solvent molecules**.** Right side bottom, illustration of the $\alpha_2\beta_2$-**P** (blue — above and red — below the plane) with corresponding positions.

Positions	Calculated			Exp.
	Conf. A	Conf. A + 1 CH_3CN	Conf. A + 2 CH_3CN	
1 and 14	1,7	0,0	0,0	0,98
2 and 13	11,5	0,0	0,0	0,59
2^1 and 13^1	1,2	0,0	0,0	0,26
2^2 and 13^2	2,4	0,0	0,0	0,22
3 and 12	6,7	0,0	0,0	0,22
3^1 and 12^1	3,2	0,0	0,0	0,12
3^2 and 12^2	6,7	0,0	0,0	0,12
4 and 11	1,2	0,0	0,0	0,40
6 and 9	1,3	0,0	0,0	0,35
7 and 8	3,0	0,0	0,0	0,88
7^1 and 8^1	1,1	0,0	0,0	0,19
7^2 and 8^2	4,1	0,0	0,0	0,02
10 and 5	2,6	0,0	0,0	0,47
10^1 and 5^1	1,2	0,0	0,0	0,06
10^2 and 5^2	0,4	0,0	0,0	0,09
10^3 and 5^3	2,6	0,0	0,0	0,14
10^4 and 5^4	1,1	0,0	0,0	0,00
10^5 and 5^5	0,3	0,0	0,0	0,00
10^6 and 5^6	0,2	0,0	0,0	0,04
15 and 20	2,0	0,0	0,0	0,19
15^1 and 20^1	1,5	0,0	0,0	0,00
15^2 and 20^2	0,2	0,0	0,0	0,30
15^3 and 20^3	0,2	0,0	0,0	0,00
15^4 and 20^4	0,6	0,0	0,0	0,00
15^5 and 20^5	2,7	0,0	0,0	0,00
15^6 and 20^6	3,7	0,0	0,0	1,30
16 and 19	2,9	0,0	0,0	0,52
17 and 18	3,8	0,0	0,0	0,20

0-1.40 ppm	0.0-0.14 ppm
1.50-2.9 ppm	0.15-0.30 ppm
3.0-6.0 ppm	0.31 -0.4 ppm
> 6.0 ppm	> 0.4 ppm



17^1 and 18^1	0,5	0,0	0,0	0,12
17^2 and 18^2	2,2	0,0	0,0	1,18
21 and 23	7,0	-12,2	-13,7	1,67
22 and 24	7,4	-11,8	-10,4	0,00

NBO partial charges calculations and analysis

In order to clarify the source of Δσ in ^{13}C and ^{15}N NMR, the NBO partial charges were calculated for the following complexes: two porphyrin cations 4+ (differing by the position of protons), porphyrin cation 6+, optimized $[\alpha_2\beta_2$-**P**$[H_2SO_4]_2]^{4+}$ and $[\alpha_2\beta_2$-**P**$[H_2SO_4 \cdot H_2O]_2]^{4+}$ complexes (corresponding to the crystal structure and achiral environment), and conformer **A** of $\alpha_2\beta_2$-**P·10CSA(R)** without inclusion of solvent molecules and with addition of one and two acetonitrile molecules (Figure S15 and Table S9). The presence or absence of protons change the N(21) – N(24) NBO partial charges in the porphyrin cations 4+ by 0.019-0.043, wherein in the porphyrin cation 6+ where all N(21) – N(24) are protonated the partial charges are very similar and differ only by 0.001. In the achiral $[\alpha_2\beta_2$-**P**$[H_2SO_4]_2]^{4+}$ complex one of oxygen atoms of each H_2SO_4 is situated symmetrically in respect to the core protons that result in the similar partial charges of N(21) – N(24). However, addition of two water molecules leads to the difference by 0.008-0.015 in the N(21) – N(24) NBO partial charges. This increased difference in partial charges can be explained by the non-symmetric position of oxygens of H_2SO_4 with respect to the porphyrin's core protons and as a result of the H-bonds' formation of different strength. That is in good agreement with the observation that protonation or deprotonation has a stronger influence on the N(21) – N(24) NBO partial charges, than the H-bonds formation. In the more asymmetric $\alpha_2\beta_2$-**P·10CSA(R)** complex, the NBO partial charges split up to 0.056 for N(21) – N(24) following the above noticed trend. Addition of acetonitrile molecules results in decrease of the difference between the NBO partial charges of N(21)-N(24), herein the difference between the NBO partial charges of N(22) and N(24) is much more sensitive to the presence of solvent molecules in the cavity. This agrees well with our NMR calculations.

The *ortho*-carbons in the porphyrin cations 4+ and 6+ have similar partial charges (differing up to 0.006) (Figure S15 and Table S9). However, in the $[\alpha_2\beta_2$-**P**$[H_2SO_4]_2]^{4+}$ complex, due to the small size of the guest molecule only one H-bond (instead of two) can be formed between the NH_3^+ groups of porphyrin and oxygens of each H_2SO_4 guest molecule. This leads to differences between the NBO partial charges of *ortho*-carbons of up to 0.021 for phenyls with the H-bonded and non-bonded NH_3^+ groups. Addition of two water molecules allows the H-bond formation with all four NH_3^+ groups of porphyrin, although these bonds are not equal; this is proved by difference in the NBO partial charges of *ortho*-carbons of up to 0.12 between phenyls with the H-bonded NH_3^+ groups. In $\alpha_2\beta_2$-**P·10CSA(R)**, the larger size of camphorsulfonic acid allows the H-bond formation with all four NH_3^+ groups of porphyrin; however non-equality of the formed bonds results in differences between the NBO partial charges of *ortho*-carbons of up to 0.13. Wherein, the tilt of phenyl rings varied less than 5.0˚ in the porphyrin cation 6+,

$[\alpha_2\beta_2$-**P**$[H_2SO_4]_2]^{4+}$, $[\alpha_2\beta_2$-**P**$[H_2SO_4 \cdot H_2O]_2]^{4+}$ and $\alpha_2\beta_2$-**P·10CSA(R)** complexes, which allowed us to exclude this factor from consideration. Inclusion of acetonitrile molecules into the model also had a minor effect on *ortho*-carbons.

Summarizing, based on the six model complexes and their NBO partial charges it was shown that Δσ values rely on the interactions with the chiral guests of certain size and chirality transfer effect.

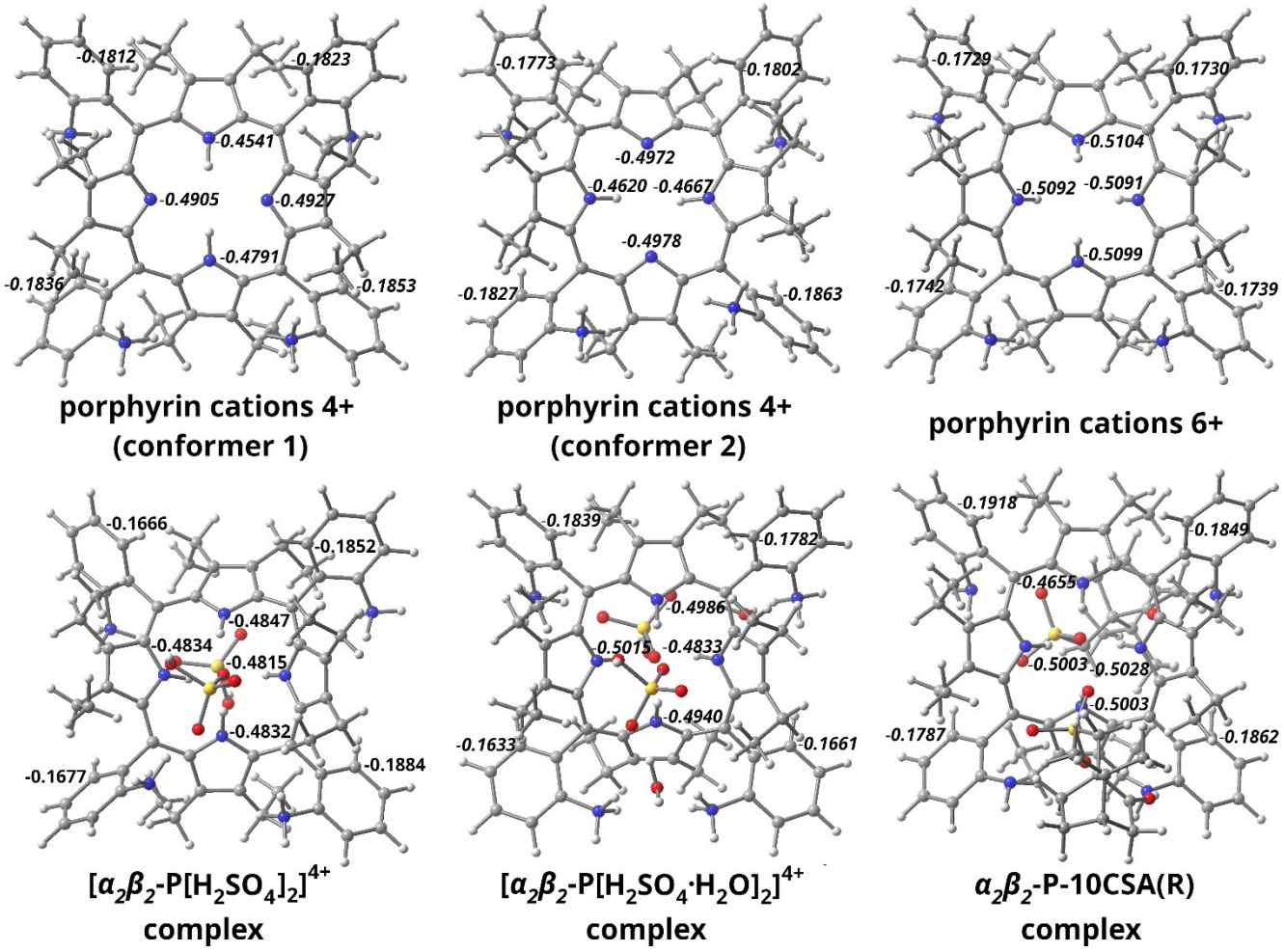


Figure S15. NBO partial charges in variety of calculated complexes.

Table S9: Partial charges for core N atoms and *ortho*-carbons in different porphyrin complexes.

Atom positions	**$[\alpha_2\beta_2$-P$]^{4+}$ conf. 1**	**$[\alpha_2\beta_2$-P$]^{4+}$ conf. 2**	**$[\alpha_2\beta_2$-P$]^{6+}$**	**$\alpha_2\beta_2$-P$[H_2SO_4]_2]^{4+}$**	**$\alpha_2\beta_2$-P$[H_2SO_4 \cdot H_2O]_2]^{4+}$**	**$\alpha_2\beta_2$-P·10CSA(R)**	**$\alpha_2\beta_2$-P·10CSA(R) + 1 CH_3CN**	**$\alpha_2\beta_2$-P·10CSA(R) + 2 CH_3CN**
N(22)	-0.47909	-0.49783	-0.50992	-0.4832	-0.49395	-0.50025	-0.49708	-0.49635
N(23)	-0.49272	-0.46670	-0.50905	-0.4815	-0.48329	-0.50283	-0.49102	-0.49120
N(24)	-0.45405	-0.49720	-0.51043	-0.4847	-0.49861	-0.46549	-0.49102	-0.50509
N(21)	-0.49051	-0.46203	-0.50923	-0.4934	-0.50147	-0.52167	-0.50010	-0.49921
$C(5^6)$	-0.18363	-0.18273	-0.17424	-0.1677	-0.16333	-0.17866	-0.17550	-0.17476
$C(10^6)$	-0.18532	-0.18629	-0.17385	-0.1884	-0.16609	-0.18624	-0.18431	-0.18450
$C(15^6)$	-0.18225	-0.18021	-0.17303	-0.1852	-0.17821	-0.18491	-0.18454	-0.18664
$C(20^6)$	-0.18124	-0.17726	-0.17285	-0.1666	-0.18393	-0.19176	-0.19280	-0.19389

Geometries

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Conformer **A**

C	2.75945500	1.52518200	2.09538300
C	-0.82229900	-3.66595300	0.98781700
C	0.13152300	-3.55406300	2.01257500
C	0.97671200	-2.41616700	1.68910400
C	-0.57190400	-2.58876600	0.04321200
C	-1.20924400	-2.33173900	-1.19933100
C	0.21127600	4.35461400	1.36816700
C	-0.87750600	4.24667800	0.48976300
C	-0.77574100	2.94962000	-0.16584600
C	-1.40952700	2.51428000	-1.35673300
C	-1.32782500	1.19718100	-1.88417800
C	-1.54333400	0.71584600	-3.24495200
C	-1.46252800	-0.68573200	-3.19904700
C	-1.21227500	-1.05512400	-1.81823200
C	4.11574700	-0.29060300	2.53089300
C	2.74912600	-0.74379000	2.30880700
C	2.20733600	-2.04960700	2.31363100
C	2.23652300	2.81515500	1.85409900
N	1.98710900	0.38066100	2.06528700
H	0.96827900	0.37495700	1.97543800
C	4.12689800	1.10208500	2.40249300
C	5.34459000	-1.13481900	2.73792800
H	5.19344000	-1.84769600	3.56898000
H	6.17011300	-0.47335500	3.06457700
N	0.26145900	2.25515000	0.44418700
H	0.73851700	1.43526300	-0.01499700
N	-1.06563400	0.10491600	-1.09061200
H	-1.28930700	0.11148900	-0.06139700
N	0.48121300	-1.83235100	0.53880100
H	1.03912500	-1.16023800	-0.04261900
C	5.79442700	-1.90359700	1.48641900
H	6.04199000	-1.21050100	0.65920600
H	4.99982800	-2.58368900	1.12494400
H	6.69325400	-2.51275100	1.70493400
C	2.96219300	-3.15598000	2.95382900
C	3.31559000	-4.31232400	2.21791500
C	4.01217900	-5.37623300	2.79951100
C	4.01283000	-4.17791000	4.91714800
C	3.32720100	-3.11255500	4.31686200
H	3.04734900	-2.22943600	4.91037400
N	2.90773500	-4.42533400	0.80135500
H	3.10552200	-3.54296100	0.17013800
H	1.87615200	-4.57451300	0.74168000
H	3.36715500	-5.24156400	0.36268500
C	0.11025900	-4.30094300	3.32133400
H	1.09049200	-4.76993700	3.53749000
H	-0.60691600	-5.14109000	3.23905900
O	1.77718700	2.19396300	-2.28218500
C	-0.29138900	-3.39458800	4.49838100
H	0.43596400	-2.57146500	4.64882500
H	-0.34015900	-3.97796900	5.43867700
H	-1.28558000	-2.93346400	4.33007100
C	-2.02693900	-4.56377900	1.01384400
H	-2.19059200	-5.03693100	0.02829400
H	-1.84564500	-5.39840100	1.71891500
O	3.78244700	1.66819700	-0.89176100
C	-3.28992600	-3.79638000	1.44076100
H	-3.19527500	-3.42316300	2.48084900
H	-4.18592200	-4.44508700	1.39036400
H	-3.46412000	-2.91662700	0.79186500
C	-1.87893300	-3.47008600	-1.88444900
O	1.71018800	0.25874500	-0.69556000
C	-1.11391500	-4.59715500	-2.26409600
H	-0.03459500	-4.58913500	-2.05288700
N	-4.13202400	-2.36927800	-1.80709300
H	-4.87698100	-2.59220000	-1.07305700
H	-3.68108800	-1.51228100	-1.35973800
C	-1.38850700	-1.60890600	-4.38391200
H	-1.77736400	-1.08651000	-5.27996700
C	0.05357600	-2.07879600	-4.65188600
H	0.70832000	-1.21680100	-4.89312600
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N	-4.31064900	2.52149700	-1.61484800
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H	-1.57506900	6.26735900	0.38997800
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H	-2.49737900	5.33812000	2.55995400
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H	1.46633200	5.88404600	2.23424700
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H	0.79215400	4.25549500	4.12081000
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H	0.36752500	5.96173100	4.46355600
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C	3.63432500	4.69054800	0.94661600
C	4.46100200	5.81152200	1.09159800
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C	4.77410700	6.27394400	2.37973800
H	5.41656800	7.15881200	2.49643300
C	4.27131000	5.60567000	3.50858100
H	4.51923800	5.96436700	4.51824500
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H	3.07674500	3.94040400	4.23327600
N	3.32405600	4.19849900	-0.40489100
H	3.82665000	4.74668600	-1.12262000
H	2.31270000	4.26355300	-0.63331600
H	3.57192900	3.12147300	-0.57829100
C	6.20690800	2.02517100	1.21774500
H	7.11442600	2.63588900	1.39191500
H	5.63377800	2.48072900	0.38977600
H	6.53726600	1.02596900	0.87029800
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H	6.00840200	1.48493800	3.30736500
H	5.14541000	2.94750400	2.86774400
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C	0.94982200	3.10350000	1.29767600
C	-3.26387600	-3.49297500	-2.16954500
C	-3.86277900	-4.59679300	-2.79182100
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H	-3.55137500	-6.55934600	-3.66069300
C	-1.69799600	-5.69528400	-2.90676000
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C	-1.38181400	4.53381200	-2.85085200
C	-4.17396900	4.45850500	-3.14528700
C	-2.02565300	5.47755900	-3.66355200
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H	-5.27005900	4.43883400	-3.25053600
H	-1.43618300	6.24623400	-4.18433800
H	-3.93465000	6.17917600	-4.44388500
H	-2.25853000	2.41601900	-4.37295200
H	-2.09563600	0.92283000	-5.29786900
C	-0.21714000	1.97868700	-4.98425200
H	0.40480400	1.10490400	-5.26697900
H	-0.30489400	2.63088900	-5.87531300
H	0.33146400	2.52296900	-4.19194600
H	-2.55556700	5.13559500	-0.53693000
S	-3.64090700	0.43175500	0.83469300
O	-2.26399200	0.02661900	1.29510200
O	-5.96296800	-2.33404700	0.10513600
H	-2.03642200	-2.49591600	-4.24257000
S	2.56756200	1.10444900	-1.62910600
C	4.06135300	-2.17071500	-1.68019100
C	4.29981000	-0.98709500	-2.60476500
C	5.97805600	-2.49281800	-3.07035500
C	5.16018800	-3.18276200	-1.96215700
H	5.74048100	-3.41322800	-1.04369600
H	4.69633700	-4.13386700	-2.29878200
C	4.86628600	-1.75124700	-3.88689400
C	6.74650000	-1.33835900	-2.39195300
H	7.48618000	-0.88262700	-3.07637300
H	7.31046700	-1.68897500	-1.50513900
H	6.60337400	-3.18541100	-3.66390900
C	5.41879400	-0.81564400	-4.96942100
H	4.60262600	-0.28812200	-5.50401600
H	6.12365900	-0.05163200	-4.59215100
H	5.96012800	-1.41278900	-5.73145400
C	3.82380200	-2.67857100	-4.53435700
H	3.07792800	-2.09246200	-5.10796000
H	4.31031600	-3.36703100	-5.25433300
H	3.25899500	-3.30289100	-3.81208300
C	3.16586200	-0.02522800	-2.92076000
H	2.24504000	-0.54593900	-3.25221000
H	3.47993600	0.66655100	-3.72812300
C	5.61547000	-0.33992000	-2.02416700
H	5.74539800	0.65014300	-2.50212300
H	5.51638900	-0.13398500	-0.94302700
O	3.15593400	-2.31216700	-0.84249000
C	4.36272400	-5.30691900	4.15835100
H	4.27594900	-4.12763100	5.98383900
H	4.28012400	-6.26512400	2.20704500
H	4.90535100	-6.14398800	4.62111800
C	-6.68197400	-1.60361600	0.78769300
C	-6.24649800	-0.40864200	1.63803300
C	-8.18273800	-1.74917600	0.96805100
C	-7.23695100	-0.57683600	2.86966700
C	-6.86416900	0.83014200	0.89590400
C	-8.50872900	-0.60819300	1.95313300
H	-8.70719900	-1.68551600	-0.00853300
H	-8.39945700	-2.76317000	1.36424500
C	-7.20488800	0.60134600	3.85089600
C	-6.98318200	-1.87266400	3.65933900
H	-6.42730400	1.75353400	1.32284300
H	-6.59905400	0.81237300	-0.18058400
C	-8.39438900	0.71210200	1.15755000
H	-9.46408200	-0.73864400	2.49535000
H	-8.77145400	1.57487500	1.73826400
H	-8.98260900	0.67586700	0.22026900
C	-4.77560400	-0.35660000	2.00824400
H	-4.34729100	-1.36627300	2.15733100
H	-4.60679600	0.23223500	2.93301500
O	-3.88528000	1.93339000	0.89463900
H	-7.31011300	1.59367200	3.37395400
H	-6.26488600	0.61054200	4.44084900
H	-8.03347800	0.49846500	4.58122900
H	-6.05529000	-1.79182400	4.26234600
H	-6.88281600	-2.77907200	3.02878600
H	-7.81191100	-2.05882700	4.37208900

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Conformer **B**

C	-6.25658600	-1.79258000	-1.34605000
C	-5.84619100	-0.66271400	-2.28257800
C	-8.12674400	-0.54648400	-2.14369900
C	-7.76448100	-1.70914800	-1.19233200
H	-8.23724700	-2.68268700	-1.43794900
H	-8.01808900	-1.50286100	-0.13099600
C	-6.92248400	0.43956200	-1.94125100
C	-7.92090700	-1.06813500	-3.58165400
H	-8.30706700	-0.35442500	-4.33352800
H	-8.44536800	-2.02845200	-3.75077200
H	-9.11985300	-0.09965200	-1.94877000
C	-6.90532100	1.63299200	-2.90372500
H	-5.98244900	2.23257900	-2.76355800
H	-6.96500900	1.35113400	-3.97147400
H	-7.76515900	2.30191200	-2.69401800
C	-6.85033600	0.98421800	-0.50003000
H	-6.10678900	1.80081300	-0.45555400
H	-7.83321500	1.40439300	-0.20511100
H	-6.56855800	0.21627200	0.24851300
C	-4.36033000	-0.37393600	-2.39646500
H	-4.14726900	0.37230800	-3.18824100
C	-6.37672200	-1.21178600	-3.67341200
H	-5.94172500	-0.59406800	-4.48451200
H	-6.03392500	-2.25135000	-3.84802800
O	-5.50643700	-2.64742600	-0.87302800
C	2.94090500	-0.27851200	-2.23029900
C	-1.18758700	4.13720600	-0.03326100
C	-0.18369000	4.39839400	-0.97862700
C	0.69527800	3.24116100	-1.00957500
C	-0.91554200	2.81617800	0.52836800
C	-1.50140900	2.19077000	1.66067000
C	0.60076700	-3.31986500	-2.35451000
C	-0.38760900	-3.61960800	-1.40329800
C	-0.27403800	-2.63777400	-0.33227600
C	-1.00957000	-2.57737000	0.88556800
C	-1.15209400	-1.39955100	1.66325200
C	-1.57255500	-1.22785000	3.04535000
C	-1.68008800	0.15319900	3.27510500
C	-1.35544600	0.82344000	2.01747700
C	4.10305100	1.70449900	-2.07672900
C	2.67866400	1.95323600	-1.87543800
C	1.98791900	3.15683800	-1.60579800
C	2.54052000	-1.62522200	-2.39595800
N	2.03975900	0.73345900	-1.96401300
H	1.02313100	0.61608700	-1.98680100
C	4.25997000	0.33051800	-2.30119100
C	5.25807500	2.66269100	-1.94649900
H	4.92971900	3.70046300	-2.11120200
H	5.97260200	2.44982700	-2.76975900
N	0.73988200	-1.75559900	-0.68310900
H	1.20203500	-1.08579500	-0.02225700
N	-0.97718900	-0.14597400	1.11702500
H	-1.11010900	-0.00955600	0.08261900
N	0.16612300	2.28598000	-0.15898200
H	0.74041800	1.47603500	0.20792900
C	5.99405500	2.57667100	-0.59616800
H	6.43538500	1.57356500	-0.42977700
H	5.30717100	2.78623900	0.24525900
H	6.82526800	3.30866100	-0.57001100
C	2.69051800	4.45119400	-1.81165100
C	2.99149400	5.27964500	-0.70422800
C	3.67364900	6.49111000	-0.86231400
C	3.74747200	6.11010200	-3.26560300
C	3.08358400	4.88748500	-3.09430300
H	2.86012100	4.25128800	-3.96317500
N	2.56769700	4.86560300	0.64499600
H	2.78307800	3.78976500	0.87187700
H	1.54160600	4.99230700	0.75919600
H	3.03319100	5.43944400	1.36811600
C	-0.15707500	5.57801400	-1.91573600
H	0.81179000	6.11220500	-1.86315500
H	-0.91594400	6.31137900	-1.58038800

O	3.18301600	2.39731300	1.03254200
C	-0.43394100	5.17647900	-3.37560200
H	0.33580700	4.47858600	-3.76138500
H	-0.43561200	6.07116600	-4.02874100
H	-1.41696600	4.67459800	-3.47800400
C	-2.37143700	5.04077300	0.19492600
H	-2.87238500	4.79916600	1.14602400
H	-2.00106700	6.08088000	0.31708400
O	1.92504300	1.50126400	3.02536600
C	-3.40307800	4.99877400	-0.95081000
H	-2.98605800	5.38668700	-1.90045800
H	-4.27589200	5.63115300	-0.69374700
H	-3.75078800	3.96558500	-1.13972400
C	-2.32260600	3.03946000	2.57062700
O	1.79181700	0.33087200	0.80156300
C	-1.71505700	3.99907600	3.40331800
H	-0.62190900	4.11229600	3.36732400
N	-4.37426900	1.93759500	1.73033300
H	-5.40395400	1.97967200	1.80688700
H	-4.11500100	2.04607300	0.66335800
C	-1.87583100	0.78621800	4.62459100
H	-2.35392100	0.04890100	5.29964800
C	-0.54343900	1.25402300	5.24067200
H	0.13236400	0.39513800	5.42421300
H	-0.72528300	1.75957100	6.20941100
H	-0.00014500	1.94835000	4.57184400
C	-1.69892400	-2.29975300	4.09402200
N	-3.96107100	-2.91745300	1.22986600
H	-3.55292800	-1.99439300	0.87602400
H	-4.50659500	-2.67723100	2.07544300
H	-4.67119800	-3.12299700	0.46135600
C	-1.50148100	-4.60632500	-1.61413800
H	-1.18026500	-5.35913700	-2.36042400
C	-2.78241400	-3.91289900	-2.10671700
H	-2.62145200	-3.43191900	-3.09315000
H	-3.11724100	-3.12550900	-1.40525300
H	-3.60946100	-4.64118800	-2.21583100
C	0.69815900	-3.90697800	-3.74034100
H	0.04929300	-4.80283300	-3.79410700
H	1.72174700	-4.26987200	-3.96018400
C	0.27609100	-2.90084500	-4.82571000
H	0.93899200	-2.01241400	-4.84052600
H	-0.75954500	-2.54200200	-4.66004200
H	0.31851500	-3.37028000	-5.82798900
C	3.42165300	-2.54890400	-3.15080400
C	3.84923600	-3.78420700	-2.60598600
C	4.68108900	-4.65660400	-3.31454400
H	5.00698500	-5.60881500	-2.86730500
C	5.09488200	-4.31412200	-4.61285300
H	5.74430600	-5.00175400	-5.17383300
C	4.67149900	-3.10457500	-5.18700800
H	4.98275300	-2.83931200	-6.20789500
C	3.85217800	-2.23158300	-4.45865200
H	3.51732200	-1.28471700	-4.90752600
N	3.38279700	-4.19059400	-1.26551900
H	3.90341700	-5.02338700	-0.94210800
H	2.37010900	-4.44101900	-1.30003600
H	3.45991300	-3.40528900	-0.49315800
C	6.02234100	-1.19945400	-1.24282000
H	7.01786300	-1.65029100	-1.42252800
H	5.30666800	-2.01684200	-1.03635900
H	6.09304100	-0.58322800	-0.32421200
C	5.59053300	-0.35815400	-2.45393100
H	6.35781100	0.41728600	-2.64430300
H	5.59261200	-0.99839400	-3.35591900
O	-3.75358900	1.75240100	-0.81618900
C	1.33276100	-2.16623700	-1.86509600
C	-3.72542300	2.89359400	2.63119900
C	-4.50602000	3.67384000	3.49070200
H	-5.60051000	3.55597300	3.51684700
C	-3.87901500	4.61674000	4.32325300
H	-4.48895400	5.22764800	5.00444100
C	-2.48492600	4.77868900	4.27933300
H	-1.99535300	5.51667300	4.93146300
H	-4.09046100	0.94249500	1.87774600
C	-1.63332900	-3.84247900	1.36500000
C	-3.02926000	-4.01537100	1.50584800
C	-0.80534000	-4.94510900	1.67713700
C	-3.57796600	-5.23832000	1.91521100
C	-1.34063000	-6.16418700	2.11050900
H	0.28338100	-4.82489800	1.57559300
C	-2.73277800	-6.31518600	2.22134300
H	-4.66967100	-5.35352900	1.99759500
H	-0.67415500	-7.00306000	2.35762700
H	-3.16591800	-7.27198300	2.54719700
H	-2.31592700	-3.14279700	3.72732000
H	-2.24427000	-1.88432500	4.96412800
C	-0.34312200	-2.84927300	4.56712300
H	0.26743400	-2.05226000	5.03455100
H	-0.48895000	-3.65166800	5.31665400
H	0.23383900	-3.27613200	3.72190400
H	-1.71340300	-5.17676900	-0.69247400
S	-3.42108800	0.27437300	-0.97890500
O	-3.82754500	-0.45650300	0.30713000
O	-1.98130700	0.03417300	-1.36043100
H	-2.57774700	1.63853000	4.57149300
S	2.63609400	1.16993400	1.76682400
C	3.81003200	-2.32182200	1.73318600
C	3.98809100	-1.16575400	2.69361500
C	4.56628200	-3.05288800	3.87184800
C	4.17160000	-3.59332100	2.48114600
H	3.32089300	-4.30801900	2.49460900
H	5.00895300	-4.10220400	1.95862800
C	5.24836300	-1.68850100	3.51571100
C	3.25455100	-2.61681500	4.56004400
H	3.42026300	-2.35829000	5.62257300
H	2.50020300	-3.42743400	4.54660500
H	5.19291000	-3.74767100	4.46121900
C	5.57971800	-0.82222400	4.73674200
H	6.04137600	0.13973400	4.43362500
H	4.70905100	-0.58715700	5.37665800
H	6.32537800	-1.34442000	5.37044400
C	6.51323400	-1.80917300	2.64872500
H	6.90584200	-0.80607900	2.38416300
H	7.31679800	-2.33002700	3.20737600
H	6.36603000	-2.36018200	1.69772200
C	4.14688000	0.21591500	2.08512100
H	4.65741900	0.18655700	1.10309200
H	4.71923900	0.88201100	2.76136000
C	2.82062300	-1.37518900	3.73082900
H	2.72922600	-0.45423200	4.33760800
H	1.84799300	-1.49255700	3.21369300
O	3.45013300	-2.24554100	0.54772200
C	4.05034700	6.90819900	-2.15000800
H	4.03781100	6.43870900	-4.27419900
H	3.91170600	7.11806600	0.01119500
H	4.58168900	7.86280100	-2.27585100
H	-3.80639000	-1.30076700	-2.64658000

200

Conformer **C**

C	-2.74727900	1.60768600	-2.03318300
C	0.69464400	-3.69427300	-0.99021800
C	-0.25179800	-3.53994400	-2.01574800
C	-1.06739500	-2.38432000	-1.67703500
C	0.47050200	-2.62270600	-0.03248200
C	1.12973900	-2.37636500	1.20021500
C	-0.12351600	4.37822400	-1.34704600
C	0.97734900	4.25178900	-0.48607300
C	0.86125000	2.95686300	0.17274800
C	1.53724400	2.48814800	1.32383600
C	1.45106000	1.16206800	1.82941100
C	1.67143300	0.67809500	3.18732800
C	1.50080400	-0.71747200	3.16076000
C	1.21433900	-1.08862800	1.78945800
C	-4.15257800	-0.16862600	-2.47877300
C	-2.79726900	-0.65889200	-2.26889200
C	-2.28859400	-1.97788800	-2.29426000
C	-2.19003700	2.88354900	-1.79487600
N	-2.00526700	0.44257700	-2.01491300
H	-0.98603600	0.40799400	-1.93955300
C	-4.12739200	1.22231800	-2.33497300
C	-5.40317800	-0.97973000	-2.68730900
H	-5.27368100	-1.68900700	-3.52505700
H	-6.21331100	-0.29499500	-3.00427200
N	-0.20581000	2.28452200	-0.41119400
H	-0.68852100	1.47300700	0.05826300
N	1.13683800	0.07146700	1.04900900
H	1.25656300	0.03062500	0.01357300
N	-0.55944000	-1.83220900	-0.51767300
H	-1.09423400	-1.14704700	0.07120500
C	-5.86607800	-1.74739000	-1.43998100
H	-6.08854700	-1.05571400	-0.60443500
H	-5.08837000	-2.45283600	-1.09069700
H	-6.78281700	-2.32937700	-1.65795200
C	-3.07344300	-3.05650100	-2.94551400
C	-3.45468000	-4.21045100	-2.21994900
C	-4.17905900	-5.25088000	-2.80985200
C	-4.15496400	-4.03126500	-4.91546200
C	-3.44115200	-2.98945200	-4.30678000
H	-3.14119500	-2.10755400	-4.89219900
N	-3.04777000	-4.34524800	-0.80478900
H	-3.21292000	-3.45980200	-0.16995500
H	-2.02185400	-4.53085800	-0.74936500
H	-3.53490700	-5.14631500	-0.36813500
C	-0.24391100	-4.25813300	-3.34030700
H	-1.23242800	-4.70482000	-3.56688900
H	0.45834100	-5.11236900	-3.27726900
O	-1.70506200	2.22198400	2.35210600
C	0.17469200	-3.33088900	-4.49518200
H	-0.54465500	-2.49862000	-4.63355100
H	0.22542100	-3.89461900	-5.44726900
H	1.17023800	-2.88057300	-4.30810300
C	1.87328600	-4.62655400	-1.02881500
H	2.02011700	-5.12179500	-0.05193300
H	1.66480300	-5.44382100	-1.74672000
O	-3.70328400	1.73730000	0.93400500
C	3.16769600	-3.90495700	-1.44259100
H	3.08610200	-3.50169800	-2.47150200
H	4.02609200	-4.60390300	-1.40965100
H	3.37634300	-3.04487900	-0.77670600
C	1.76572600	-3.52828800	1.88612600
O	-1.64790800	0.29728100	0.75689300
C	1.00175200	-4.68084000	2.18050000
H	-0.06858300	-4.67205400	1.92765400
N	3.97529400	-2.35925200	2.05288000
H	4.97816900	-2.61536500	2.06715800
H	3.83910400	-1.73191100	1.15057000
C	1.36751900	-1.61482400	4.36449800
H	1.76111700	-1.08833100	5.25600000
C	-0.09654100	-2.02056200	4.61588700
H	-0.71916400	-1.12532900	4.81701100
H	-0.17453700	-2.69085700	5.49413900
H	-0.52391100	-2.55262000	3.74125300
C	1.81386600	1.49986300	4.43713600
N	4.40838600	2.29584000	1.30961800
H	4.36543800	1.35869200	1.75335200
H	5.40854600	2.45202800	1.06368200
H	3.93700000	2.09539500	0.33614500
C	2.12865300	5.21788200	-0.43756000
H	1.74771100	6.23548900	-0.65702600
C	3.23628200	4.85888900	-1.44879400
H	2.85774300	4.88564000	-2.49002700
H	3.63905900	3.84162800	-1.27882800
H	4.06828800	5.58682600	-1.37415000
C	-0.33285000	5.50545000	-2.32023600
H	0.34050400	6.33962500	-2.04261000
H	-1.35860200	5.91285700	-2.23465400
C	-0.05688700	5.09434900	-3.77767100
H	-0.70021000	4.25099600	-4.09860400
H	0.99513200	4.77404900	-3.91484000
H	-0.24464900	5.94513100	-4.46194500
C	-3.04540500	4.09198500	-1.99059500
C	-3.53177200	4.78320200	-0.85811500
C	-4.33415400	5.92399400	-0.98394600
H	-4.71265100	6.44386600	-0.08986400
C	-4.65289800	6.40305100	-2.26460000
H	-5.27627500	7.30326300	-2.36650800
C	-4.17962200	5.73242100	-3.40494600
H	-4.43120400	6.10541500	-4.40849200
C	-3.39102300	4.58153000	-3.26674600
H	-3.03044000	4.04714800	-4.15746600
N	-3.21527800	4.27261100	0.48556300
H	-3.70326400	4.81953200	1.21449100
H	-2.20114600	4.32190800	0.70568200
H	-3.47585900	3.19807500	0.64689800
C	-6.16555800	2.19305900	-1.11765500
H	-7.05645300	2.83204200	-1.27518300
H	-5.56728500	2.62560500	-0.29507800
H	-6.52109100	1.20182000	-0.77226700
C	-5.35881400	2.08440000	-2.42302500
H	-6.01140100	1.65644500	-3.21168700
H	-5.10516500	3.09461100	-2.77961400
O	4.07344400	-0.50727000	0.30423100
C	-0.88949000	3.14395600	-1.25676200
C	3.13631900	-3.55212200	2.23790200
C	3.72616100	-4.68180200	2.81763400
H	4.79808400	-4.68179300	3.07068300
C	2.94525400	-5.81723200	3.08607400
H	3.41128600	-6.70230300	3.54282800
C	1.57481100	-5.80940000	2.78048000
H	0.95591500	-6.68986000	3.00599000
H	3.82048800	-1.68676200	2.83062800
C	2.36946400	3.45618600	2.09452200
C	3.77885100	3.35215300	2.09629700
C	1.77204900	4.50024300	2.83008400
C	4.57406000	4.26458400	2.80020700
C	2.55964600	5.40653700	3.55356900
H	0.67626500	4.58618800	2.83306600
C	3.95998200	5.29158500	3.53584200
H	5.67114300	4.17991300	2.77340900
H	2.07954400	6.21132800	4.12930700
H	4.58305600	6.00644600	4.09285700
H	2.50111200	2.34992500	4.27353200
H	2.28751900	0.87802200	5.22345000
C	0.45534500	2.02202900	4.94277000
H	-0.20011600	1.18415800	5.25637600
H	0.59756600	2.68441000	5.81905600
H	-0.08879400	2.57751400	4.15519500
H	2.55940800	5.27678100	0.57561700
S	3.50872400	0.02269200	-1.02707200
O	2.22229000	-0.65450100	-1.41656500
O	6.21255500	1.54077400	-0.44929300
H	1.97522600	-2.53659400	4.26280800
S	-2.50671000	1.15150200	1.68417000
C	-4.11100000	-2.07488900	1.70451200
C	-4.30447500	-0.89655500	2.64538200
C	-6.02285800	-2.35634400	3.10963500
C	-5.23583600	-3.05798300	1.98640800
H	-5.83137400	-3.26100200	1.07132200
H	-4.79721200	-4.02632800	2.30803400
C	-4.88208200	-1.65793600	3.92395400
C	-6.76213500	-1.17138900	2.45147200
H	-7.48090000	-0.70129900	3.14832900
H	-7.34515500	-1.49458200	1.56657900
H	-6.66358800	-3.03653500	3.70108900
C	-5.39586900	-0.71869400	5.02234100
H	-4.55886600	-0.22535400	5.55729900
H	-6.07759500	0.07316300	4.66021500
H	-5.95120100	-1.30733100	5.78091000
C	-3.86215200	-2.62372900	4.55113800
H	-3.09417500	-2.06720000	5.12510700
H	-4.36281900	-3.30580300	5.26746900
H	-3.32293500	-3.25612200	3.81648400
C	-3.13935800	0.02613600	2.96335900
H	-2.23453300	-0.52585600	3.28737200
H	-3.42844100	0.72054100	3.77776300
C	-5.60476100	-0.20356200	2.08340400
H	-5.70075100	0.78489700	2.57266400
H	-5.50837200	0.01022200	1.00351300
O	-3.22015300	-2.23053500	0.85367300
C	-4.53085700	-5.15890900	-4.16704000
H	-4.41982600	-3.96368500	-5.98073300
H	-4.46792600	-6.13872300	-2.22572000
H	-5.09552600	-5.97730300	-4.63679400
C	6.77788700	0.60256800	-0.99704300
C	6.20426500	-0.30830600	-2.08806700
C	8.20895000	0.13695100	-0.75769500

C	6.94321900	-1.66992800	-1.76557200
C	6.99001900	0.17947100	-3.37140500
C	8.36625600	-1.01079600	-1.77553700
H	8.91877100	0.97877200	-0.89532300
H	8.31734500	-0.18594100	0.29890300
C	6.74792700	-2.74601800	-2.84130600
C	6.56964800	-2.29032300	-0.40708300
H	6.50460200	-0.24961100	-4.27094500
H	6.92792800	1.28151600	-3.47394300
C	8.43563400	-0.35406600	-3.17250900
H	9.19647100	-1.70647000	-1.54916400
H	8.71134600	-1.08522200	-3.95590300
H	9.19312200	0.45284500	-3.20676400
C	4.70152400	-0.32524800	-2.36035800
H	4.38858200	-1.31626300	-2.74478800
H	4.40640700	0.43138900	-3.11487000
O	3.32448300	1.52630500	-0.93741700
H	6.94651600	-2.39658300	-3.87131800
H	5.71393900	-3.14935300	-2.82218700
H	7.42654100	-3.60220700	-2.64915900
H	5.55246200	-2.72766700	-0.45699200
H	6.58977100	-1.56123100	0.42678000
H	7.27304200	-3.11257000	-0.16310400

200

Conformer **D**

C	-5.06953200	-1.06405600	-1.15437300
C	-4.55250800	-1.16068600	-2.57317600
C	-6.81462500	-1.52338700	-2.71224400
C	-6.56484100	-1.31494000	-1.20078000
H	-6.86165200	-2.17629100	-0.56556300
H	-7.09012800	-0.42659300	-0.78892800
C	-5.78613900	-0.53588500	-3.36324200
C	-6.23520800	-2.90742900	-3.07836400
H	-6.51050600	-3.20162900	-4.10836600
H	-6.61737700	-3.70304000	-2.40965700
H	-7.86637900	-1.36477300	-3.01532400
C	-5.66904400	-0.66774400	-4.88698900
H	-4.89569500	0.01826900	-5.28891200
H	-5.42598500	-1.68693000	-5.23968400
H	-6.62969700	-0.37619000	-5.35894100
C	-6.02553300	0.94416800	-3.02801000
H	-5.25629500	1.58208100	-3.50899300
H	-7.00782800	1.27676900	-3.41975800
H	-6.00912200	1.16889600	-1.94340200
C	-3.21117500	-0.50421800	-2.81746600
H	-3.14962300	0.48235800	-2.32109800
C	-4.70267000	-2.68899500	-2.93132500
H	-4.14363700	-2.87808200	-3.86972300
H	-4.23483500	-3.32380000	-2.15795700
O	-4.38262200	-0.78729100	-0.15718500
C	2.48452900	1.26611000	-2.03842600
C	-3.15311800	2.67721900	0.88589200
C	-2.57515100	3.45596400	-0.12736900
C	-1.26364900	2.88696400	-0.42163000
C	-2.22545400	1.57798800	1.16185700
C	-2.31680900	0.54571800	2.13281900
C	2.25534600	-2.50624800	-2.67865200
C	1.58337200	-3.42573300	-1.86113000
C	1.14139000	-2.69879100	-0.66728700
C	0.40927100	-3.18679400	0.44475100
C	-0.34608200	-2.40188900	1.36483200
C	-0.77460500	-2.75493500	2.71319700
C	-1.49017400	-1.65800000	3.22083900
C	-1.56556700	-0.66025700	2.15264400
C	2.33535300	3.53920000	-1.71850400
C	1.02482300	2.91610200	-1.51112800
C	-0.18508500	3.52625500	-1.09905800
C	2.92689500	-0.03675100	-2.36863500
N	1.17286400	1.55911200	-1.71422300
H	0.38797100	0.86882400	-1.76733500
C	3.23577100	2.51453800	-2.01739600
C	2.75398900	4.97597400	-1.54268000
H	1.93168600	5.66256400	-1.79637800
H	3.54277200	5.18617600	-2.29438700
N	1.56283200	-1.38815800	-0.80506100
H	1.54264500	-0.66007100	-0.05575100
N	-0.83682000	-1.14441100	1.08976000
H	-0.88609100	-0.76398500	0.10475000
N	-1.14476100	1.73324900	0.31893700
H	-0.22546400	1.24278300	0.47536300
C	3.29717000	5.32023900	-0.14496700
H	4.19252800	4.71695200	0.10620000
H	2.55166900	5.12428100	0.64701500
H	3.58990500	6.38781200	-0.10249600
C	-0.28483000	5.01065300	-1.16518300
C	-0.29940400	5.76746000	0.02987400
C	-0.36278000	7.16548600	0.00761700
C	-0.44815900	7.09997900	-2.42481700
C	-0.35869900	5.70120900	-2.39217600
H	-0.34048100	5.12357000	-3.32793300
N	-0.27662100	5.07205800	1.32979900
H	0.45070100	4.22175300	1.39641100
H	-1.21073000	4.66064500	1.53503200
H	-0.05523500	5.73234000	2.09375700
C	-3.28116600	4.54209300	-0.89823200
H	-2.71947500	5.49599300	-0.87577200
H	-4.24860000	4.75765900	-0.40442400
O	1.47136800	3.17359700	1.41871100
C	-3.53520500	4.13785400	-2.36042200
H	-2.58939200	3.93282400	-2.90076000
H	-4.06536400	4.94575300	-2.90192200
H	-4.16013400	3.22372500	-2.41824700
C	-4.52638000	2.94083300	1.45007000
H	-4.56798800	2.69238500	2.52340300
H	-4.70936200	4.03310600	1.39979000
O	0.72028300	1.63347900	3.26669100
C	-5.66519700	2.22236000	0.71077100
H	-5.70494600	2.52276300	-0.35535500
H	-6.64183600	2.47433700	1.16844000
H	-5.53952600	1.12318800	0.74008800
C	-3.28538200	0.78338500	3.24050500
O	1.28479600	0.71055000	0.99428800
C	-3.02008800	1.78021000	4.20144900
H	-2.07260900	2.33445500	4.12893400
N	-4.72823600	-1.06981300	2.43295400
H	-4.67957200	-0.84280700	1.34904200
H	-4.00430700	-1.80168300	2.58300200
C	-1.83468900	-1.47608500	4.67720200
H	-1.88329900	-2.47110900	5.16231500
C	-0.78829900	-0.60358200	5.39901400
H	0.19837400	-1.10815000	5.41727900
H	-1.09522000	-0.42069000	6.44750800
H	-0.64457000	0.37031600	4.89281500
C	-0.36597400	-3.96085200	3.51702900
N	-2.02486000	-4.67807800	0.02205700
H	-1.90684700	-3.74726400	-0.58228000
H	-2.53414900	-4.37231500	0.87089500
H	-2.66476700	-5.29871600	-0.50292500
C	1.25173100	-4.83144400	-2.28638400
H	2.03526600	-5.16969300	-2.99386800
C	-0.11921400	-4.96432700	-2.97307600
H	-0.14953600	-4.41521100	-3.93269100
H	-0.92957200	-4.52776400	-2.35971300
H	-0.34640200	-6.03069700	-3.16941300
C	2.65645700	-2.73340400	-4.11477600
H	2.56127400	-3.81384900	-4.33875500
H	3.72504400	-2.49032600	-4.28429400
C	1.78754800	-1.93150000	-5.10108700
H	1.90962200	-0.83920200	-4.95679900
H	0.71379900	-2.15926100	-4.96267800
H	2.07637500	-2.16755000	-6.14416300
C	4.16301800	-0.20191400	-3.17135300
C	5.21196100	-1.05830200	-2.75639000
C	6.37916700	-1.21787400	-3.51003400
H	7.18520300	-1.88293000	-3.16195100
C	6.51481400	-0.52939700	-4.72710500
H	7.42929500	-0.65856600	-5.32404500
C	5.48099700	0.30782500	-5.17711200
H	5.57641400	0.83634000	-6.13684300
C	4.32582400	0.47359200	-4.40206600
H	3.51504700	1.12914900	-4.75288400
N	5.06459800	-1.84934000	-1.51828700
H	5.95898700	-2.30873200	-1.27724600
H	4.34832700	-2.59667100	-1.65871100
H	4.71652100	-1.28122900	-0.64232200
C	5.49822400	2.15873300	-0.89753200
H	6.58726800	2.31256800	-1.02920400
H	5.31439600	1.07879400	-0.74296800
H	5.19673700	2.68091300	0.03354100
C	4.72805500	2.69600600	-2.11472800
H	4.94089400	3.77683100	-2.22660500
H	5.12649800	2.22354800	-3.03066900
O	-1.04408500	-2.00607800	-3.39106900
C	2.24942200	-1.22989000	-1.98547700
C	-4.48615800	0.05408400	3.35640000
C	-5.42016100	0.32462300	4.36239900
H	-6.35610900	-0.25154700	4.43296100
C	-5.14392900	1.33492100	5.29937500
H	-5.87383300	1.55074700	6.09304300
C	-3.93643400	2.04915500	5.22887600
H	-3.71128400	2.82473800	5.97554400
H	-5.64782100	-1.50176500	2.62614600
C	0.38495800	-4.66530200	0.62984000
C	-0.79476700	-5.39950700	0.39166800
C	1.54314100	-5.38097100	0.99780600
C	-0.83287700	-6.79407100	0.50626200
C	1.51687800	-6.77522800	1.13845100
H	2.47569400	-4.82423600	1.16877800
C	0.32998900	-7.48372400	0.88583400
H	-1.76097100	-7.35022200	0.30157800
H	2.42815700	-7.31571900	1.43340400
H	0.30473300	-8.57908400	0.97962100
H	-0.47312100	-4.88859000	2.92401800
H	-1.06386600	-4.07238800	4.37026600
C	1.07065900	-3.86706100	4.05597600
H	1.17989600	-3.01131400	4.75044300
H	1.34377100	-4.78888000	4.60604400
H	1.80061000	-3.73200900	3.23299600
H	1.31544900	-5.52828600	-1.43409000
S	-1.72795700	-1.41124000	-2.22188800
O	-2.21618000	-2.43010900	-1.18484500
O	-0.91036300	-0.33424600	-1.49509700
H	-2.83506200	-1.02823000	4.81887300
S	1.56336000	1.78287200	2.05077500
C	4.34943300	-0.60390000	1.76448600
C	3.85139700	0.30030800	2.87340600
C	5.24562000	-1.21986900	3.88491000
C	5.25532100	-1.64412500	2.40059200
H	4.87549500	-2.67235500	2.21819100
H	6.26244500	-1.58968200	1.93641000
C	5.15645600	0.34053800	3.78471800
C	3.86017200	-1.61032300	4.44356900
H	3.81466400	-1.48374700	5.54141900
H	3.62794500	-2.67282700	4.23844300
H	6.10541800	-1.60204900	4.46593700
C	4.93788500	1.03770900	5.13257500
H	4.85130400	2.13648200	5.00776500
H	4.04205400	0.69311600	5.68182500
H	5.81575700	0.86220200	5.78752400
C	6.34671400	1.01078600	3.07688800
H	6.19538700	2.10730300	3.00716900
H	7.28089600	0.85175300	3.65235800
H	6.52994300	0.64704300	2.04506300
C	3.32683100	1.65988600	2.45027500
H	3.83083500	2.03323300	1.53800600
H	3.45279000	2.40785200	3.25834200
C	2.89771900	-0.62950500	3.71418000
H	2.31661700	0.00809800	4.40754300
H	2.15166900	-1.12813100	3.06327300
O	4.09008200	-0.49253400	0.55689100
C	-0.43909300	7.83294400	-1.22612800
H	-0.51379900	7.62244700	-3.39036800
H	-0.35939700	7.74148500	0.94622800
H	-0.49409900	8.93114400	-1.24483600
H	-3.01346300	-0.37997100	-3.89987100

200

Conformer **E**

C	3.10275700	1.45297900	-3.08939600
C	3.65925200	0.11472500	-3.62537900
C	5.01021900	1.73059400	-4.52831100
C	4.02141700	2.53214500	-3.66136000
H	3.42928900	3.28999400	-4.21411900
H	4.51273900	3.06815000	-2.82471500
C	5.20981000	0.42353800	-3.68826200
C	4.22359700	1.24031500	-5.76316200
H	4.89013300	0.78891700	-6.52213700
H	3.68423100	2.06893900	-6.26178200
H	5.94696500	2.27133000	-4.76198400
C	6.02306700	-0.65778100	-4.41240000
H	6.11166600	-1.57714700	-3.79619600
H	5.61365400	-0.95478800	-5.39624600
H	7.05782800	-0.29690800	-4.58415900
C	5.86113400	0.67361300	-2.31876500
H	5.98824000	-0.27741000	-1.76577700
H	6.86919600	1.11673200	-2.45209100
H	5.27262300	1.34164500	-1.66224000
C	3.21016000	-1.21925500	-3.02976400
H	4.00127200	-1.98748400	-3.14572200
C	3.25865600	0.18647500	-5.15301300
H	3.39148500	-0.81548400	-5.60920500
H	2.18819500	0.44926000	-5.26632100
O	2.10808400	1.64732900	-2.40419000
C	-0.56994600	3.01514300	-0.49052300
C	-1.17697900	-2.98203900	-2.56631200
C	-1.56649500	-1.84327000	-3.28616300
C	-1.67888600	-0.75113800	-2.33461600
C	-1.05419200	-2.58186100	-1.16612200
C	-0.67327000	-3.36458600	-0.04797800
C	2.82412900	2.90116600	1.36983700
C	3.37633600	1.77176300	1.98555000
C	2.38855400	0.69310500	1.87504700
C	2.38631100	-0.57739900	2.49671300
C	1.41620400	-1.60185900	2.29208800
C	1.09575100	-2.76658000	3.11694700
C	0.15114200	-3.52806600	2.40565100
C	-0.07127500	-2.85521100	1.13589600
C	-2.57294200	2.93564500	-1.62709300
C	-1.82044600	1.70802000	-1.85627300
C	-2.19970700	0.55526400	-2.58036400
C	0.50605900	3.39108700	0.34109300
N	-0.62474300	1.81452800	-1.17292700
H	0.20432300	1.21964800	-1.35453900
C	-1.81176900	3.73993700	-0.77711300
C	-3.96672100	3.26881600	-2.08876600
H	-4.05868700	3.12725000	-3.18115700
H	-4.13699900	4.34958800	-1.91822000
N	1.33335600	1.18490100	1.12535500
H	0.38036500	0.74647000	1.06380500
N	0.63843700	-1.67837800	1.16243100
H	0.85467800	-1.16579800	0.26395700
N	-1.34723800	-1.23444800	-1.09283200
H	-1.48893100	-0.69280100	-0.20464000
C	-5.07774300	2.47931200	-1.38152800
H	-5.06440700	2.66021000	-0.28943700
H	-4.96097500	1.39081000	-1.53909100
H	-6.07256700	2.77638000	-1.76734000
C	-3.25155400	0.64375600	-3.62111800
C	-4.39539400	-0.19094700	-3.57220900
C	-5.39255200	-0.13577500	-4.55096200
C	-4.12738500	1.59376100	-5.70408500
C	-3.14230400	1.53797000	-4.70865800
H	-2.25473900	2.18497400	-4.77005000
N	-4.53926700	-1.17265700	-2.47529700
H	-4.34754400	-0.79644100	-1.45711600
H	-3.85882400	-1.95328400	-2.60514500
H	-5.48880700	-1.58274300	-2.48706900
C	-1.59479700	-1.70970900	-4.78678500
H	-2.54875500	-1.26676200	-5.13437600
H	-1.55416500	-2.71949200	-5.24021700
O	-0.53017600	0.44862700	3.47778100
C	-0.42367200	-0.86195200	-5.31265700
H	-0.42755600	0.15660300	-4.87386200
H	-0.47294500	-0.76286500	-6.41469200
H	0.54826400	-1.33125700	-5.05948200
C	-0.72291700	-4.26751300	-3.20258600

H	-0.97753400	-5.13972000	-2.57735400
H	-1.27735400	-4.41357600	-4.15116500
O	-1.93922000	2.30084500	2.53758300
C	0.78800600	-4.25889600	-3.49905100
H	1.04198100	-3.49596300	-4.26260600
H	1.11879100	-5.24088300	-3.89054200
H	1.38434900	-4.02194300	-2.59491800
C	-0.76341200	-4.84264300	-0.16733300
O	-1.27709200	0.37392200	1.08462500
C	-1.99741400	-5.50229300	-0.34031800
H	-2.91821100	-4.90158700	-0.35113500
N	1.71090900	-4.98158800	0.09312700
H	1.92195800	-4.06554700	-0.44797900
H	1.77363400	-4.70462000	1.09844000
C	-0.67884500	-4.65796800	2.95791400
H	-0.24210400	-4.99234100	3.91920500
C	-2.14164600	-4.22893500	3.17908100
H	-2.20032800	-3.38004000	3.89066900
H	-2.73606500	-5.06389200	3.59918600
H	-2.62110900	-3.91693600	2.22813500
C	1.47356100	-2.99446000	4.55105900
N	4.63104900	-2.02718800	1.47251800
H	3.95352600	-2.78566400	1.26172700
H	5.56262200	-2.37182300	1.18586000
H	4.32246700	-1.26053600	0.73483400
C	4.78204900	1.74195800	2.52475100
H	5.01107100	2.74493300	2.93837100
C	5.84036400	1.39996200	1.45987100
H	5.85299300	2.14814600	0.64325400
H	5.65321000	0.41904700	0.98436200
H	6.84986500	1.37906700	1.91494500
C	3.55222000	4.18943100	1.09915800
H	4.47635600	4.21702200	1.70837200
H	2.95694700	5.06252700	1.42666400
C	3.91486300	4.33989200	-0.38879300
H	3.02328500	4.24503400	-1.04010200
H	4.63089000	3.55065500	-0.69544000
H	4.38801600	5.32278800	-0.58126600
C	0.56518800	4.78520500	0.86500000
C	0.33347500	5.02356000	2.23931400
C	0.37176600	6.31825700	2.77188700
H	0.17768900	6.48899600	3.84245400
C	0.66698700	7.40258900	1.92990100
H	0.70441600	8.41868900	2.34890200
C	0.91423400	7.18687100	0.56352900
H	1.14896500	8.03554100	-0.09536900
C	0.85292900	5.88962900	0.03637400
H	1.03073700	5.71589600	-1.03483600
N	0.04230800	3.89105400	3.13597200
H	-0.18336900	4.22975500	4.08661000
H	0.86393600	3.26204500	3.23162200
H	-0.80732600	3.22273600	2.82778900
C	-3.21511600	4.93808400	0.99918800
H	-3.49338700	5.94552500	1.36631100
H	-2.72880000	4.38194300	1.82165100
H	-4.15508700	4.40340000	0.75664600
C	-2.29900200	5.05664600	-0.23080000
H	-2.84961700	5.57447100	-1.04338600
H	-1.45508800	5.72067400	0.00788600
O	3.75945400	-0.49028000	-0.46376100
C	1.51737700	2.52116300	0.85512500
C	0.40798400	-5.63203500	-0.15281300
C	0.36855600	-7.01925400	-0.32418700
H	1.29463600	-7.61522200	-0.32069700
C	-0.87385600	-7.65250900	-0.49788100
H	-0.91129700	-8.74329800	-0.63180500
C	-2.05675700	-6.89481500	-0.49391800
H	-3.02942000	-7.39258600	-0.61911100
H	2.48656000	-5.63568500	-0.10980300
C	3.53219800	-0.90442800	3.39859400
C	4.63709700	-1.61506900	2.88628000
C	3.57618700	-0.48555100	4.74320100
C	5.76070500	-1.89686700	3.67399700
C	4.68837200	-0.77394100	5.54730700
H	2.72383400	0.07602000	5.15092800
C	5.78237500	-1.47525000	5.01269500
H	6.62048100	-2.44093300	3.25233400
H	4.70597900	-0.44173100	6.59559300
H	6.66107100	-1.69563900	5.63601900
H	2.54601800	-2.78152700	4.71343200
H	1.33970400	-4.06707500	4.79657300
C	0.61557300	-2.13533800	5.50142900
H	-0.44594700	-2.45378200	5.47211800
H	0.97001400	-2.23794000	6.54599400
H	0.63326700	-1.06507200	5.21689500
H	4.87440300	1.05121500	3.37746900
S	2.75522100	-1.28973300	-1.27781000
O	2.82770600	-2.76046300	-0.83737100
O	1.34531500	-0.75595600	-1.19428500
H	-0.65874600	-5.54455100	2.29298000
S	-1.60748800	0.80809700	2.51036800
C	-4.74786800	-0.21466400	0.95516400
C	-4.44367400	0.21000200	2.38327900
C	-6.73848700	0.19546200	2.20853000
C	-6.26063200	-0.23298900	0.80822300
H	-6.59419700	0.43587900	-0.01283200
H	-6.58395800	-1.26055600	0.53809000
C	-5.67142500	-0.47342100	3.13956500
C	-6.42180800	1.70201200	2.33580300
H	-6.87421500	2.13884200	3.24561600
H	-6.82191500	2.27676500	1.47721200
H	-7.78587400	-0.08193900	2.43014700
C	-5.79645400	-0.06225900	4.61184100
H	-5.01044500	-0.53766100	5.23346900
H	-5.74676600	1.02902800	4.78395100
H	-6.76959900	-0.41360900	5.01164800
C	-5.62770000	-2.00959700	3.06724000
H	-4.84676800	-2.41408500	3.74317700
H	-6.59261900	-2.44266600	3.39963700
H	-5.41855100	-2.41221200	2.05495600
C	-3.09490300	-0.11390800	3.00138500
H	-2.79590100	-1.16956300	2.84100000
H	-3.13479600	0.06428200	4.09500300
C	-4.86990000	1.72566000	2.40131900
H	-4.48516600	2.17709900	3.33572200
H	-4.38341200	2.28598700	1.58203000
O	-3.93645200	-0.49476800	0.05887500
C	-5.25628800	0.76182000	-5.62349300
H	-4.01461200	2.28990900	-6.54800200
H	-6.27674400	-0.78978000	-4.49136300
H	-6.03615800	0.80287700	-6.39770800
H	2.28829100	-1.60408400	-3.50961000

Conformer **F**

C	-5.03020800	-2.08569500	-3.78559700
C	-4.85881500	-0.91341000	-2.78052400
C	-7.10866200	-1.16660400	-3.06162300
C	-6.53119700	-2.23279400	-4.01464900
H	-6.78276700	-2.07029800	-5.08323100
H	-6.83823800	-3.27240200	-3.77811800
C	-6.10185700	-1.18555200	-1.85132600
C	-6.83589000	0.21493900	-3.70373400
H	-7.36900900	1.02574000	-3.17114300
H	-7.17920900	0.24848400	-4.75580300
H	-8.16301600	-1.34009700	-2.77335800
C	-6.41525800	-0.08784800	-0.82089800
H	-5.75994000	-0.21256300	0.06362900
H	-6.28398100	0.93740000	-1.21323400
H	-7.46566900	-0.18928600	-0.47850800
C	-6.03319000	-2.53026800	-1.11068200
H	-5.26191700	-2.47840300	-0.31420400
H	-7.00369300	-2.75930800	-0.62543400
H	-5.77585200	-3.39120500	-1.75795500
C	-3.41784700	-0.93196200	-2.31294200
H	-2.75004700	-0.56055700	-3.11602700
C	-5.29095100	0.36114800	-3.57588400
H	-4.99181400	1.26516600	-3.00919200
H	-4.77335200	0.41117100	-4.55540000
O	-4.12605400	-2.76533500	-4.23804500
C	2.40928600	-0.20055900	-2.36151300
C	-1.58541800	4.12039300	0.22766700
C	-0.70181000	4.37542900	-0.83047700
C	0.25658800	3.28400000	-0.86859200
C	-1.16434900	2.85796700	0.84450300
C	-1.64983400	2.22413200	2.01549600
C	0.29579800	-3.37551000	-1.89113500
C	-0.41888700	-3.69752600	-0.72713200
C	-0.13334800	-2.65359300	0.25751000
C	-0.65261000	-2.50957200	1.57334500
C	-0.78699400	-1.27972500	2.27400000
C	-1.17036400	-1.05157200	3.66276100
C	-1.49473300	0.31014300	3.78150100
C	-1.29638500	0.91854200	2.46671600
C	3.50630400	1.82040400	-2.42425300
C	2.12717700	2.01883300	-1.98336300
C	1.45391100	3.21223000	-1.64148500
C	2.06343200	-1.57159900	-2.41838700
N	1.52444800	0.77603500	-1.94936400
H	0.52245400	0.60947800	-1.76753500
C	3.67119200	0.45389700	-2.67575200
C	4.63233000	2.81928900	-2.48036300
H	4.24689300	3.84480400	-2.58573700
H	5.21463400	2.62982800	-3.40685000
N	0.73433700	-1.75428800	-0.33429700
H	1.24893700	-0.99051100	0.16929800
N	-0.79866400	-0.05921200	1.63797500
H	-0.86483700	0.00640700	0.59237700
N	-0.08815700	2.38466500	0.11372800
H	0.56233200	1.60942100	0.41730900
C	5.57288500	2.76238600	-1.26181200
H	6.07283500	1.77733100	-1.17203900
H	5.01607300	2.95115400	-0.32424300
H	6.36994500	3.52543500	-1.36180400
C	2.05213200	4.52338400	-2.00296700
C	2.44096100	5.42941000	-0.98711200
C	3.01026100	6.67023500	-1.29465100
C	2.79263500	6.15865900	-3.66515300
C	2.24359600	4.90935800	-3.34690400
H	1.95126600	4.21188900	-4.14532700
N	2.24416800	5.06391800	0.42775900
H	2.54746700	4.01607400	0.67007000
H	1.24436200	5.15543700	0.69681800
H	2.78545700	5.69278000	1.04509100
C	-0.87270800	5.45290400	-1.86872400
H	0.05200400	6.04966800	-1.99202200
H	-1.64174800	6.16770200	-1.51708800
O	3.06621300	2.65901800	0.86690500
C	-1.29471700	4.87541900	-3.23193300
H	-0.51589200	4.20719400	-3.65064400
H	-1.46800700	5.69011500	-3.96214800
H	-2.22826300	4.28321700	-3.14745400
C	-2.77325300	5.00061400	0.52185300
H	-3.04371000	4.97116700	1.58917800
H	-2.46652900	6.04952400	0.33135900
O	2.15207600	1.65712900	3.00016400
C	-4.01255900	4.69868600	-0.34057000
H	-3.79584400	4.79673800	-1.42238700
H	-4.82883400	5.40482700	-0.09210000
H	-4.39340800	3.67106100	-0.19096400
C	-2.68941800	2.95623400	2.80191400
O	1.77969300	0.51894000	0.78904400
C	-2.36070700	3.96904400	3.72413700
H	-1.30147600	4.20519500	3.89835000
N	-4.42302900	1.66329900	1.58030200
H	-5.44980100	1.56043100	1.50739600
H	-4.05464100	1.82555600	0.56846500
C	-1.73594000	1.03072900	5.07522600
H	-2.08545700	0.30773700	5.83855600
C	-0.45096900	1.72373200	5.57165500
H	0.33431300	0.97977600	5.81200100
H	-0.65600700	2.31298400	6.48708100
H	-0.02192500	2.39727300	4.80357100
C	-1.05614600	-2.02433000	4.80927300
N	-3.47801900	-2.63593000	2.19023700
H	-3.25055100	-2.05772700	1.30408900
H	-3.38545600	-1.94434600	2.96457400
H	-4.47344900	-2.91365500	2.12958000
C	-1.43261400	-4.80800400	-0.64649800
H	-1.13832400	-5.60495300	-1.35800300
C	-2.85942300	-4.34484300	-0.98582600
H	-2.95321100	-4.04938500	-2.05063100
H	-3.15543100	-3.47026300	-0.37283200
H	-3.59028600	-5.15650600	-0.80151100
C	0.12795900	-4.02293100	-3.24278800
H	-0.45456800	-4.95728200	-3.12286500
H	1.10575900	-4.33669500	-3.65974000
C	-0.57942400	-3.10683100	-4.25892100
H	-0.04004300	-2.14618600	-4.38568000
H	-1.62356800	-2.88248000	-3.96235900
H	-0.62333100	-3.59970400	-5.24994200
C	2.82057000	-2.46893100	-3.32287200
C	3.43813900	-3.65136500	-2.84667100
C	4.15416400	-4.50333100	-3.69276000
H	4.63348000	-5.41355400	-3.29947000
C	4.25450900	-4.19272900	-5.05967300
H	4.81237900	-4.86455600	-5.72806600
C	3.63986000	-3.03482700	-5.56331000
H	3.70799400	-2.79531100	-6.63450300
C	2.93913500	-2.18082700	-4.70086000
H	2.45421700	-1.27487800	-5.09376600
N	3.30794500	-4.01764700	-1.42184300
H	3.95957300	-4.78548600	-1.18676200
H	2.33956200	-4.35070600	-1.22270800
H	3.47678400	-3.18508800	-0.71576900
C	5.68673900	-0.97555400	-1.96664100
H	6.66682900	-1.34342300	-2.32833200
H	5.09384700	-1.85046600	-1.64257100
H	5.86973000	-0.35106300	-1.06951800
C	4.97630800	-0.18292800	-3.07546000
H	5.65173900	0.62232300	-3.42485500
H	4.83065700	-0.84188900	-3.95201300
O	-3.49470300	1.41282000	-0.90486000
C	1.04827700	-2.16435300	-1.61303100
C	-4.05371000	2.67257800	2.58383700
C	-5.06671200	3.36931100	3.25387100
H	-6.12692100	3.14284400	3.06114400
C	-4.71733600	4.36832100	4.17635800
H	-5.50913900	4.91586000	4.70780100
C	-3.36470500	4.66603900	4.41218900
H	-3.09029200	5.45109300	5.13186500
H	-4.02519200	0.71072300	1.74696400
C	-1.23575100	-3.71923900	2.21462700
C	-2.62265700	-3.80805000	2.46181300
C	-0.43853800	-4.82937600	2.56192700
C	-3.20946800	-4.95472200	3.00659000
C	-1.00614400	-5.97946800	3.12899800
H	0.64444000	-4.77978300	2.37686100
C	-2.39327400	-6.04847000	3.33964900
H	-4.29562000	-5.00466600	3.18077900
H	-0.36678700	-6.83221300	3.39943800
H	-2.84754400	-6.95300200	3.76940100
H	-1.65555100	-2.93972800	4.62946700
H	-1.49441600	-1.55709000	5.71243600
C	0.39483100	-2.43482600	5.10628000
H	1.01436300	-1.55248100	5.36057500
H	0.43161100	-3.13740400	5.96164600
H	0.85637100	-2.93953000	4.23434600
H	-1.42088600	-5.28670000	0.34716400
S	-2.89886800	0.02197500	-0.88414100
O	-3.36705600	-0.71081500	0.39387900
O	-1.38444700	0.02148600	-0.98792300
H	-2.55033200	1.77089600	4.97082800
S	2.70149900	1.39558000	1.64875100
C	4.09598000	-2.02921100	1.40539400
C	4.30455500	-0.86304200	2.34681700
C	5.14971900	-2.70527500	3.43223000
C	4.63960300	-3.27077100	2.09013300
H	3.85525900	-4.05110800	2.19198400
H	5.44853500	-3.70800200	1.46758500
C	5.68386400	-1.29298400	3.01478000
C	3.89496900	-2.37023700	4.26798200
H	4.15843200	-2.10383600	5.30855700
H	3.20589600	-3.23535700	4.32959600
H	5.88879700	-3.35240200	3.93976500
C	6.08868400	-0.40769100	4.19936700
H	6.43769400	0.58783800	3.85640300
H	5.28302100	-0.24132200	4.93843800

H	6.94039300	-0.87301000	4.73626200
C	6.84460000	-1.31683600	2.00568100
H	7.13043900	-0.28639600	1.71109800
H	7.74341200	-1.78134300	2.45896500
H	6.62612800	-1.87104400	1.07008000
C	4.29363600	0.52696400	1.73389300
H	4.65940900	0.52914400	0.68878600
H	4.91865500	1.22403700	2.32726600
C	3.28007700	-1.16187000	3.50712000
H	3.18926400	-0.25187600	4.13002300
H	2.26597000	-1.34987500	3.10206200
O	3.58004200	-1.98156500	0.27770800
C	3.18185100	7.03704400	-2.63998500
H	2.92421000	6.44672400	-4.71834400
H	3.32057500	7.35830200	-0.49270100
H	3.62309900	8.01483200	-2.88221200
H	-3.07157200	-1.95312500	-2.06147300

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Conformer **G**

C	4.97627900	-3.55935300	-2.25116000
C	4.76542200	-2.02772900	-2.14822000
C	6.67413200	-2.31037600	-3.36805400
C	6.24721000	-3.76326800	-3.06790800
H	6.99214100	-4.35396200	-2.49536200
H	6.00696300	-4.35891100	-3.97287900
C	5.30061900	-1.57262600	-3.55598100
C	7.20002700	-1.72081200	-2.03726500
H	7.70113300	-0.74636700	-2.19286800
H	7.94234200	-2.39021800	-1.56040500
H	7.37156200	-2.21002900	-4.22121700
C	5.42111100	-0.05188500	-3.71894800
H	4.42078100	0.42561700	-3.70013700
H	6.02648600	0.43590100	-2.93261600
H	5.88980300	0.18777200	-4.69575200
C	4.44627000	-2.11385100	-4.71522400
H	3.49308100	-1.55069000	-4.80003900
H	4.97422500	-1.98309300	-5.68164900
H	4.18317000	-3.18712900	-4.62461600
C	3.32602100	-1.74066700	-1.77989600
H	2.89161900	-2.60790900	-1.23955600
C	5.90678000	-1.56309400	-1.18155700
H	5.73053600	-0.50660200	-0.89954000
H	5.91636000	-2.15764000	-0.24706600
O	4.24510600	-4.41191800	-1.77274700
C	-1.01428300	2.04436900	-2.25927100
C	-0.22090400	-3.90601400	-0.07982800
C	-0.53741900	-3.60358400	-1.41352900
C	-1.00464400	-2.22206100	-1.43266900
C	-0.49481700	-2.71495800	0.70830900
C	-0.32359300	-2.53010600	2.11159300
C	1.98730300	3.87385800	-0.67236200
C	2.67218200	3.46349700	0.47576900
C	1.89665400	2.37261300	1.07635700
C	2.06765400	1.73145800	2.32372700
C	1.31585400	0.60807100	2.78927400
C	0.97423700	0.23852000	4.16065800
C	0.19080100	-0.92748500	4.09137600
C	0.12014500	-1.31245500	2.68680500
C	-2.60512000	0.75203500	-3.31617700
C	-1.64047300	-0.10246100	-2.62830000
C	-1.62602100	-1.51473500	-2.50399300
C	-0.23212500	3.11789800	-1.77654400
N	-0.70586500	0.71691700	-2.03395500
H	0.16342200	0.36150300	-1.56303900
C	-2.23304800	2.07612900	-3.07725400
C	-3.86020800	0.33195200	-4.03222700
H	-3.65228600	-0.48924300	-4.74092900
H	-4.20670100	1.17820000	-4.65676400
N	0.82824600	2.12548700	0.23429900
H	0.00919600	1.50484700	0.44795800
N	0.76291500	-0.33442800	1.95882700
H	1.00399100	-0.42052200	0.92915300
N	-0.92102000	-1.73266500	-0.14938900
H	-1.33293900	-0.81689300	0.16208400
C	-4.99995000	-0.08297300	-3.08980500
H	-5.29223500	0.75166600	-2.42287500
H	-4.69716100	-0.93076900	-2.44589500
H	-5.89389700	-0.39130400	-3.66664900
C	-2.37936800	-2.34782700	-3.47316100
C	-3.39295400	-3.23059100	-3.02928000
C	-4.11858100	-4.02999300	-3.91803800
C	-2.81177900	-3.11860600	-5.75786700
C	-2.10461600	-2.30989100	-4.85734400
H	-1.31144700	-1.63904400	-5.21912400
N	-3.67300400	-3.35187900	-1.58140300
H	-3.77016700	-2.41280100	-1.01949200
H	-2.88949600	-3.85864300	-1.11289300
H	-4.53571400	-3.90230600	-1.43147900
C	-0.24159000	-4.46522700	-2.61438900
H	-1.13898900	-4.56291500	-3.25732800
H	-0.01443700	-5.49272800	-2.26678700
O	-1.37585700	2.38362400	2.49829000
C	0.93221100	-3.94401700	-3.46290000
H	0.76146400	-2.89951500	-3.79606200
H	1.05810700	-4.56973100	-4.36828800
H	1.88223500	-3.98177600	-2.89491900
C	0.54141000	-5.09934200	0.42383600
H	0.05557600	-5.51024000	1.33006200
H	0.50472600	-5.91279700	-0.32718000
O	-2.76018400	2.93108300	0.46989900
C	2.01912400	-4.75546500	0.71506800
H	2.62506800	-4.66888600	-0.21040200
H	2.49852900	-5.53884600	1.33670400
H	2.11794300	-3.75767700	1.20797900
C	-0.58982700	-3.70494600	2.98461000
O	-1.51000300	0.76433800	0.59907500
C	-1.85120200	-4.33849200	2.95503500
H	-2.62307800	-3.93177400	2.28478000
N	1.71775900	-3.63031300	3.89150600
H	2.16643900	-3.72188900	2.94954300
H	1.64416400	-2.60310400	4.10836500
C	-0.67104700	-1.46032300	5.21208300
H	-0.36203000	-0.97472100	6.15820000
C	-2.16496800	-1.18469700	4.96178000
H	-2.34703900	-0.09730500	4.84641400
H	-2.77609900	-1.54609100	5.81203900
H	-2.52291500	-1.70039600	4.04673600
C	1.17186700	1.06469100	5.39752000
N	4.70175400	0.47416200	2.62019400
H	4.05915900	0.16261700	1.75716600
H	4.73487000	-0.35056600	3.24437200
H	5.65429900	0.55794300	2.22136200
C	4.03663600	3.96625600	0.85957700
H	4.15218000	4.99646600	0.46814300
C	5.17845900	3.09440200	0.30451500
H	5.21500500	3.12830900	-0.80102300
H	5.03842300	2.02726800	0.56154400
H	6.15463600	3.44225900	0.69622600
C	2.51985400	4.81721700	-1.71411500
H	3.35259600	5.40135800	-1.27600300
H	1.75645900	5.56330900	-2.00692700
C	3.02441500	4.06425700	-2.95945300
H	2.21596500	3.47334500	-3.43506900
H	3.82921800	3.35039600	-2.69638700
H	3.41730300	4.77590600	-3.71222600
C	-0.51864700	4.51185900	-2.22039900
C	-1.01205900	5.44978900	-1.28397300
C	-1.28419100	6.77478500	-1.64690500
H	-1.67960800	7.48749700	-0.90620300
C	-1.04316800	7.19186300	-2.96539400
H	-1.24813200	8.23413800	-3.25016000
C	-0.53936300	6.28179600	-3.91013800
H	-0.34277700	6.60894400	-4.94156800
C	-0.28967900	4.95328000	-3.54078200
H	0.10254200	4.23815400	-4.27852100
N	-1.25600400	5.02850700	0.10711600
H	-1.71939100	5.78428800	0.63914400
H	-0.36776000	4.81809800	0.60182600
H	-1.89320200	4.11058100	0.23177900
C	-4.03606100	3.79537400	-2.49224300
H	-4.53582300	4.70808300	-2.87188900
H	-3.54728800	4.03521900	-1.53001100
H	-4.82512900	3.04813300	-2.27606300
C	-3.03426900	3.26623700	-3.53195500
H	-3.58671700	2.97866700	-4.44882000
H	-2.36817300	4.08618100	-3.84455500
O	3.45428100	0.93243600	-1.26055100
C	0.81704600	3.01659200	-0.80688300
C	0.37227900	-4.24779500	3.86532800
C	0.12116700	-5.35406800	4.67897000
H	0.89973700	-5.75785600	5.34494900
C	-1.14926500	-5.95430800	4.63352600
H	-1.36224400	-6.82302300	5.27297500
C	-2.13271100	-5.44406300	3.77190800
H	-3.12688600	-5.91250600	3.73919100
H	2.33145500	-4.08539400	4.59093700
C	3.09899900	2.30011000	3.24191400
C	4.36455800	1.69656900	3.37976000
C	2.85480600	3.49531000	3.95362200
C	5.35677100	2.25198000	4.20184700
C	3.83078700	4.05028900	4.79011600
H	1.87720500	3.98458200	3.84071000
C	5.08554300	3.42839900	4.91464900
H	6.34515400	1.77367200	4.28588000
H	3.61691100	4.97790800	5.34092800
H	5.86168200	3.86281500	5.56132200
H	2.17979900	1.51828300	5.40410100
H	1.13638900	0.40356800	6.28681000
C	0.09848200	2.16336500	5.53245400
H	-0.90337700	1.72282300	5.70608500
H	0.32970300	2.82669300	6.38902600
H	0.01664200	2.77401900	4.61285200
H	4.13453100	4.06113300	1.95382700
S	2.94261100	-0.34694500	-0.70911200
O	3.47909400	-0.70104000	0.68810800
O	1.40320400	-0.42075500	-0.64704700
H	-0.53468800	-2.54752200	5.37849400
S	-2.23172300	1.84206200	1.40516800
C	-4.77486000	-0.64537300	0.48911100
C	-4.86008600	0.59873200	1.35926800
C	-7.03414300	-0.06831100	1.00304600
C	-6.19913000	-1.11526200	0.24260700
H	-6.42204800	-1.17762300	-0.84317000
H	-6.31552900	-2.13996800	0.65446700
C	-6.13264600	0.23116000	2.24762600
C	-6.94575500	1.23809800	0.18340800
H	-7.64251300	2.00737000	0.56543600
H	-7.21194600	1.07131700	-0.87910100
H	-8.06124300	-0.40050100	1.24326400
C	-6.64363600	1.39351400	3.10805000
H	-5.96998600	1.58953200	3.96708900
H	-6.77045400	2.34379100	2.55679400
H	-7.63158400	1.12850300	3.53702600
C	-5.87871500	-0.97625400	3.16704400
H	-5.23238300	-0.69015400	4.02201500
H	-6.83113500	-1.34635700	3.59722200
H	-5.38898700	-1.83627100	2.66514100
C	-3.66199100	1.03706700	2.18325400
H	-3.20054300	0.19545800	2.73613500
H	-3.98192300	1.79472900	2.92687500
C	-5.46373700	1.66559900	0.37058000
H	-5.35647900	2.66360200	0.83669500
H	-4.87821400	1.71323200	-0.56608500
O	-3.75302600	-1.18781900	0.03884600
C	-3.82495900	-3.97223800	-5.29093900
H	-2.57346200	-3.08134000	-6.83089800
H	-4.90773800	-4.70590700	-3.55286800
H	-4.38971700	-4.60351000	-5.99225500
H	2.69917200	-1.55318500	-2.67433700

200

Conformer **H**

C	2.83701100	-3.38087900	1.95664100
C	3.68969000	-2.49698200	2.86533800
C	4.85429700	-4.45887400	2.67207100
C	3.63035900	-4.66774100	1.75254600
H	3.01423400	-5.55763000	1.99647500
H	3.90394500	-4.75739400	0.68056200
C	5.16419000	-2.93238000	2.48573900
C	4.33943600	-4.53772600	4.12492500
H	5.17251200	-4.56706700	4.85253400
H	3.72985000	-5.44623100	4.29894400
H	5.70603400	-5.12778400	2.44450900
C	6.22539600	-2.38561700	3.45164000
H	6.37614900	-1.29687300	3.29626800
H	5.99711100	-2.54017000	4.52289300
H	7.20284600	-2.87078900	3.25194300
C	5.61822600	-2.59740000	1.05755300
H	5.82999100	-1.51684500	0.95295100
H	6.55257700	-3.15020500	0.82778500
H	4.86989800	-2.84500200	0.28396100
C	3.30514500	-1.03190200	3.00404800
H	4.13041900	-0.45630900	3.46692500
C	3.50895400	-3.23216000	4.25808700
H	3.88838800	-2.56946900	5.06171200
H	2.43881800	-3.41123700	4.47474500
O	1.70278400	-3.13875800	1.55523000
C	0.36241600	-2.02496300	-1.88454800
C	-1.83781200	0.26571800	3.71030000
C	-1.70303600	-1.10638700	3.45409900
C	-1.43111000	-1.25398000	2.02385600
C	-1.60640600	0.95965200	2.45663100
C	-1.58048800	2.36598500	2.22345800
C	3.47031100	0.10440800	-2.69033900
C	3.79498300	1.30561200	-2.05291500
C	2.59844100	1.74983400	-1.33470400
C	2.33595600	2.98887100	-0.71404700
C	1.14480400	3.30137400	0.02628300
C	0.36556500	4.52955700	-0.03843800
C	-0.81098600	4.31366100	0.70313800
C	-0.71131800	2.98392500	1.28491000
C	-1.20884700	-3.51875500	-1.09328200
C	-0.85297100	-2.56140200	-0.05189000
C	-1.38660400	-2.44999000	1.26079100
C	1.35528100	-1.39597400	-2.66746700
N	0.08684900	-1.69560900	-0.56721300
H	0.69051700	-1.10766200	0.06531300
C	-0.50198400	-3.15648100	-2.24246500
C	-2.21312600	-4.64067100	-1.03528800
H	-2.21455400	-5.10993600	-0.03848900
H	-1.87083000	-5.43921900	-1.72417400
N	1.64090600	0.75896400	-1.47589600
H	0.61238200	0.88505800	-1.25997600
N	0.46700900	2.41850300	0.83401600
H	0.83714600	1.46108200	1.11199800
N	-1.35321300	0.01684000	1.49007100
H	-1.21864400	0.27139800	0.47681600
C	-3.64772300	-4.24324600	-1.41225400
H	-3.70934000	-3.89471200	-2.46234400
H	-4.03254300	-3.42691900	-0.77076400
H	-4.32936200	-5.11059400	-1.30968800
C	-1.93747700	-3.68861700	1.87137600
C	-3.28617500	-3.81651700	2.27037400
C	-3.78698200	-5.00667400	2.80759000
C	-1.56962100	-5.99085200	2.61529300
C	-1.08453200	-4.79820700	2.06025100
H	-0.03091800	-4.68609900	1.75840900
N	-4.19007700	-2.65061400	2.17975300
H	-4.39957700	-2.27615400	1.19981800
H	-3.79057400	-1.84753600	2.71467900
H	-5.11034400	-2.87587700	2.59633100
C	-1.54790200	-2.16628600	4.51557400
H	-2.15382600	-3.06735500	4.30432200
H	-1.92887300	-1.77087100	5.47756900
O	-2.17214200	2.29235900	-2.81043200
C	-0.07390500	-2.58136700	4.66951100
H	0.34983300	-2.91058400	3.70017900
H	0.02519500	-3.40613600	5.40240900
H	0.53704800	-1.73001700	5.03289800
C	-1.90283600	0.92211200	5.06186800
H	-2.68464800	1.70533100	5.07905000
H	-2.21268600	0.17802100	5.82158800
O	-1.51988000	-0.12222400	-3.16239900
C	-0.54789600	1.52405000	5.47593900
H	0.22330700	0.74019600	5.61168800
H	-0.62889300	2.08530100	6.42724800

H	-0.14894200	2.20085300	4.68553500
C	-2.54481400	3.18599500	3.00071900
O	-0.95800300	1.03504300	-1.00281200
C	-3.90765200	2.80919300	3.04695000
H	-4.21998600	1.91814500	2.48275700
N	-0.77385700	4.77849800	3.72381800
H	-0.17431100	4.07105300	4.19890300
H	-0.42363000	4.87099200	2.73191900
C	-2.07332600	5.13597000	0.56746600
H	-1.82571800	6.06282300	0.01444700
C	-3.17179400	4.36452400	-0.18820000
H	-2.79058400	3.95324000	-1.14412000
H	-4.03111700	5.02938400	-0.40403100
H	-3.54966600	3.51691200	0.41883000
C	0.58336700	5.67164900	-0.98681700
N	4.10746900	3.73256200	1.44586500
H	3.74965400	2.67458800	1.43566500
H	3.46482000	4.23415500	2.08268000
H	5.02363300	3.71212000	1.92718500
C	5.18731600	1.86643700	-1.96775600
H	5.76874700	1.50691100	-2.83947300
C	5.91157700	1.44507700	-0.67438300
H	6.09579300	0.35413700	-0.65482500
H	5.30599600	1.66388500	0.22583800
H	6.88683900	1.96399900	-0.59167100
C	4.44114000	-0.77041800	-3.43048500
H	5.29891100	-0.15102500	-3.75883200
H	3.98599000	-1.15734000	-4.36138700
C	4.95900500	-1.93842700	-2.57200600
H	4.12612800	-2.54203900	-2.15953600
H	5.54446000	-1.56840000	-1.70922000
H	5.61006000	-2.60432200	-3.17219300
C	1.62072400	-1.88609200	-4.05306800
C	1.18535300	-1.09985500	-5.14282200
C	1.41446100	-1.48340000	-6.47000400
H	1.05745600	-0.85984600	-7.30450300
C	2.10649300	-2.67686100	-6.72779400
H	2.29563000	-2.97997700	-7.76783700
C	2.55227900	-3.47597200	-5.66133400
H	3.09559500	-4.41073500	-5.86270900
C	2.30299000	-3.08636800	-4.33816400
H	2.65004300	-3.71048300	-3.50219400
N	0.44719100	0.15071200	-4.87297000
H	0.04071000	0.53000600	-5.74525800
H	1.07707500	0.88291900	-4.49513400
H	-0.39081800	0.04671800	-4.13774600
C	-1.62244900	-3.05571600	-4.55233200
H	-1.72970300	-3.62603700	-5.49576700
H	-1.22388600	-2.05608200	-4.80080400
H	-2.63291000	-2.90006400	-4.12953300
C	-0.71295100	-3.82108500	-3.57650500
H	-1.16500200	-4.81477500	-3.38840300
H	0.25114800	-4.03155000	-4.06855900
O	3.60817200	-0.62041200	0.31345000
C	2.11118200	-0.24009700	-2.29101900
C	-2.19113000	4.34803200	3.72441500
C	-3.11210100	5.09818900	4.45903500
H	-2.79567800	5.99296500	5.01782300
C	-4.45792500	4.69416400	4.48092900
H	-5.19171900	5.27957900	5.05339200
C	-4.85230000	3.54852100	3.77203900
H	-5.90537200	3.23207200	3.78343700
H	-0.65375800	5.68281800	4.21514800
C	3.34512600	4.07533500	-0.90104600
C	4.22134000	4.42111700	0.14793500
C	3.51668600	4.72645500	-2.14199400
C	5.23256100	5.37723800	-0.01927000
C	4.50905700	5.69927300	-2.31743200
H	2.85622700	4.45476100	-2.97728700
C	5.37197000	6.02329900	-1.25644900
H	5.91684800	5.61812600	0.80937100
H	4.61961000	6.19857800	-3.29107800
H	6.16255900	6.77582200	-1.39044300
H	1.64987200	5.96193900	-1.01279900
H	0.03907300	6.56437200	-0.61893500
C	0.09089800	5.31719100	-2.40427700
H	-1.00260200	5.14138900	-2.42194400
H	0.31942500	6.13476500	-3.11604800
H	0.56363600	4.38516400	-2.77595700
H	5.18550200	2.96602200	-2.05334500
S	2.90408600	-0.06877400	1.49700700
O	3.31098200	1.35281900	1.90258100
O	1.37265000	-0.08161800	1.33525200
H	-2.47794700	5.47774500	1.54225500
S	-1.94985900	0.97096700	-2.18105600
C	-5.53601600	-0.87316800	-1.03661500
C	-4.69442900	-0.01785300	-1.99065800
C	-6.81915400	-0.00121100	-2.83092400
C	-6.95553600	-0.86567700	-1.56011000
H	-7.32016500	-1.90033000	-1.73122000
H	-7.62577800	-0.42690400	-0.79101700
C	-5.75389800	1.07361700	-2.40913200
C	-6.05105800	-0.85920800	-3.86522300
H	-6.05715300	-0.39171400	-4.86756000
H	-6.50346600	-1.86367400	-3.97870600
H	-7.77585700	0.41679700	-3.19538800
C	-5.31103500	1.98103800	-3.56377900
H	-4.48824200	2.65133200	-3.25051100
H	-4.95808200	1.42970800	-4.45497700
H	-6.16414500	2.61300400	-3.88587800
C	-6.17648100	1.96570700	-1.23011000
H	-5.38117000	2.70343600	-0.99846600
H	-7.08429300	2.54794400	-1.48696400
H	-6.39638400	1.41245400	-0.29354000
C	-3.40720800	0.35613800	-1.29326600
H	-3.01668300	-0.53329000	-0.75625600
H	-3.58159100	1.15232200	-0.54026100
C	-4.61259200	-0.91773300	-3.27497400
H	-3.84867300	-0.49281200	-3.95264700
H	-4.27405100	-1.94161700	-3.02556000
O	-5.11194400	-1.46028700	-0.03609600
C	-2.92236100	-6.10204800	2.97603400
H	-0.89086500	-6.84416100	2.75953000
H	-4.84405000	-5.08842400	3.10579500
H	-3.31295400	-7.03869000	3.39970500
H	2.40018300	-0.89907300	3.62910600

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$[\alpha_2\beta_2\text{-}P]^{4+}$ conf. 1

C	-1.30026300	2.86766700	-0.27331600
C	4.11011300	-0.35962700	1.05008400
C	3.99713900	1.02501600	1.11362300
C	2.68778900	1.34235100	0.48946800
C	2.84836800	-0.84370600	0.42863900
C	2.57361100	-2.19210600	0.03443100
C	-4.07629000	0.56432300	1.17840700
C	-4.04317500	-0.81996000	1.04669100
C	-2.75160100	-1.13394900	0.37874300
C	-2.33639500	-2.42773000	-0.07045300
C	-0.99596600	-2.79311000	-0.37142100
C	-0.48645800	-4.04287500	-0.93399000
C	0.91774000	-3.99393100	-0.86983900
C	1.28657800	-2.70043100	-0.29572000
C	0.45414800	4.14314200	-1.04216900
C	0.96102700	2.97588000	-0.32868300
C	2.28118900	2.64261900	0.06373900
C	-2.56590100	2.39771400	0.16059400
N	-0.12816400	2.21605600	0.02822500
H	-0.07288300	1.29264900	0.47130700
C	-0.94350200	4.07903700	-1.00211400
C	1.25920400	5.11245400	-1.86546000
H	2.06055500	5.58086400	-1.26486100
H	0.60305600	5.94809500	-2.17862900
N	-2.02122000	0.00074700	0.16576400
N	0.10633100	-2.02054800	-0.05511200
H	0.05900500	-1.02505800	0.20600900
N	1.99703300	0.19800200	0.17713000
C	1.86467300	4.45925600	-3.11900800
H	1.08214200	4.05357800	-3.79064600
H	2.52664800	3.60661800	-2.85346300
H	2.46957000	5.18689500	-3.69491800
C	3.30810200	3.72162100	0.01823600
C	4.46528000	3.60537700	-0.78111600
C	5.44854600	4.59483300	-0.85048000
C	4.16139800	5.89932400	0.74861600
C	3.17892500	4.89832000	0.78575600
H	2.29157000	5.01642200	1.42499000
N	4.66071400	2.34354600	-1.53427800
H	3.85756900	2.15122800	-2.16761000
H	4.70737400	1.53387900	-0.85535400
H	5.52463100	2.36736200	-2.10468100
C	4.86517800	1.93198500	1.95757600
H	5.25989700	2.79829400	1.39203400
H	5.76189100	1.36452700	2.27824700
C	4.11961300	2.44872700	3.20116200
H	3.23245900	3.05351400	2.92637300
H	4.78538800	3.08654500	3.81581300
H	3.76495200	1.61227800	3.83649700
C	5.17251200	-1.16465800	1.74996400
H	5.50273700	-2.02262100	1.13864300
H	6.07976700	-0.54198300	1.88266600
C	4.70527400	-1.64232700	3.14023400
H	4.66123900	-0.80658900	3.86648300
H	5.38300900	-2.41084400	3.56335700
H	3.65825400	-2.02917700	3.11814700
C	3.76057500	-3.09008000	-0.14489800
C	4.65214200	-2.88178000	-1.21648600
H	4.44247900	-2.06483900	-1.92157400
N	3.16703600	-4.36050700	1.89053900
H	2.17346400	-4.28770000	1.58937500
H	3.31558100	-5.28347700	2.33820800
C	1.81972200	-5.01144800	-1.52420400
H	1.29107800	-5.98462800	-1.56117900
C	2.22751300	-4.60797500	-2.95381100
H	1.34374400	-4.52573400	-3.61626600
H	2.90981000	-5.36450100	-3.38919400
H	2.74528000	-3.62917300	-2.97138400
C	-1.25264800	-5.07707900	-1.72545000
N	-2.58592300	-4.82996900	1.56614900
H	-1.61603900	-4.67062200	1.21217300
H	-2.65037300	-5.79110800	1.94789200
H	-2.75056100	-4.15818200	2.35086700
C	-4.98541300	-1.77628900	1.72710800
H	-5.93663400	-1.25659600	1.95808400
C	-4.39615300	-2.31802000	3.04446900
H	-4.35689800	-1.53539200	3.82787900
H	-3.33298000	-2.63357800	2.92115300
C	-5.01103300	1.33493000	2.08404600
H	-5.83851600	0.66434300	2.39175800
H	-5.50252000	2.18467800	1.57126000
C	-4.29115100	1.86327000	3.33787400
H	-3.47905000	2.57156600	3.07835900
H	-3.83593100	1.03670900	3.91966400
H	-5.00357500	2.39519500	3.99913700
C	-3.69000600	3.37516700	0.19335900
C	-4.86345500	3.18102600	-0.56626000
C	-5.93185400	4.08063300	-0.56361100
H	-6.82937100	3.90004300	-1.17550600
C	-5.84623200	5.22515400	0.24722800
H	-6.68080300	5.94089100	0.26430000
C	-4.70182000	5.44388700	1.03120000
H	-4.63775300	6.33479000	1.67304100
C	-3.63434800	4.53396300	0.99625200
H	-2.73456700	4.71176600	1.60330400
N	-4.97520500	1.93492600	-1.36095900
H	-5.86708700	1.89450300	-1.88558400
H	-4.90549500	1.09870900	-0.71585400
H	-4.19300200	1.84890600	-2.04239100
C	-2.46551800	4.28761000	-3.01348900
H	-3.17778300	4.95343900	-3.53928500
H	-3.00806200	3.35418300	-2.74711900
H	-1.67369400	3.99497200	-3.73130500
C	-1.88137300	4.97366700	-1.76742800
H	-1.33324000	5.88295600	-2.08333100
H	-2.70141600	5.33849600	-1.12200800
C	-2.82331800	1.04912600	0.54464800
C	4.06827100	-4.14745600	0.73426800
C	5.18967700	-4.96776400	0.59075500
H	5.39794500	-5.77956100	1.30477600
C	6.05459500	-4.73698500	-0.49226700
H	6.94007300	-5.37507800	-0.62530100
C	5.78259800	-3.69535000	-1.39300900
H	6.45813200	-3.51496800	-2.24192800
H	3.32731700	-3.61957500	2.61390500
C	-3.42078300	-3.43443200	-0.29734100
C	-3.56939200	-4.59867900	0.48283200
C	-4.37820400	-3.23210700	-1.31288300
C	-4.60282700	-5.52100900	0.30251200
C	-5.41994700	-4.14802500	-1.52788000
H	-4.29022400	-2.33611600	-1.94385200
C	-5.53811400	-5.29000600	-0.72016300
H	-4.68757800	-6.41376800	0.94162000
H	-6.14713100	-3.96945000	-2.33345400
H	-6.35540000	-6.00763700	-0.88131600
H	-2.14333300	-5.45347800	-1.18919200
H	-0.60474600	-5.96338200	-1.87067300
C	-1.69286900	-4.54603600	-3.10309600
H	-0.82039200	-4.24232300	-3.71429200
H	-2.24503900	-5.32879300	-3.65982300
H	-2.35403100	-3.66246900	-3.01158200
H	-5.26779900	-2.61185800	1.06211700
H	2.73184100	-5.19933100	-0.93065200
C	5.29066700	5.75526800	-0.07381300
H	4.04235200	6.80277300	1.36450500
H	6.33593700	4.47191300	-1.49081500
H	6.05796700	6.54193800	-0.11260000
H	-4.99197600	-3.16218400	3.44553700

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$[\alpha_2\beta_2\text{-}P]^{4+}$ conf. 2

C	-2.62587300	1.62295600	-0.29829800
C	3.63861700	1.97986200	0.95672700
C	2.77335100	3.08530600	0.91992400
C	1.51277800	2.62889800	0.34095600
C	2.90813700	0.83046300	0.43043700
C	3.37855500	-0.47883200	0.13234700
C	-3.68755400	-1.83922000	1.13556300
C	-2.88249000	-2.97574900	0.94638700
C	-1.66547900	-2.54360900	0.26442300
C	-0.57069700	-3.32426300	-0.19930100
C	0.74201800	-2.82480700	-0.46285200
C	1.85421600	-3.62509100	-1.04007000
C	3.00548900	-2.86180900	-0.85124300
C	2.54780300	-1.58724800	-0.23377600
C	-1.95075900	3.62892400	-1.15717000
C	-0.85718200	2.88274400	-0.48067300
C	0.41107800	3.40825500	-0.10686900
C	-3.46243200	0.58533800	0.20967400
N	-1.27369100	1.63710900	-0.08729800
C	-3.07653100	2.83902800	-1.02031400
C	-1.82896800	4.85103000	-2.02778700
H	-1.39921900	5.70726800	-1.47521200
H	-2.84325800	5.18160800	-2.32911000
N	-1.76315400	-1.18216700	0.10047000
N	1.18679100	-1.59537000	-0.04919100
N	1.63160600	1.26904300	0.14419500
C	-1.00291300	4.58524900	-3.29755100
H	-1.48347300	3.83016000	-3.95082500
H	0.00216300	4.17245800	-3.05287400
H	-0.85947700	5.51070100	-3.88945000
C	0.57578200	4.89226000	-0.12653500
C	1.50449200	5.53928000	-0.96706100
C	1.65762200	6.92712100	-1.01463200
C	-0.06288400	7.11170900	0.69365600
C	-0.20773100	5.71613100	0.70853000
H	-0.94264800	5.24108200	1.37460300
N	2.37272200	4.69214400	-1.81994000
H	1.80513200	4.15463100	-2.51007600
H	2.88750000	3.99236100	-1.23174900
H	3.06217800	5.25790300	-2.34738500
C	3.02794400	4.38597900	1.64731400
H	2.81359000	5.27534000	1.02529800
H	4.10800800	4.45154600	1.88487700
C	2.21229800	4.48413800	2.94990100
H	1.12257900	4.44795500	2.75390100
H	2.43255000	5.43504000	3.47417400
H	2.44950300	3.64940700	3.63930400

C	4.97782400	1.97988700	1.64628800
H	5.70415700	1.34479800	1.11218300
H	5.40854500	3.00002900	1.60900500
C	4.87531200	1.54814400	3.12323800
H	4.40524300	2.33093900	3.75060000
H	5.87241400	1.33261600	3.55612100
H	4.21023700	0.66213500	3.25305800
C	4.86831700	-0.64115200	0.08660100
C	5.62050800	-0.12443000	-0.98807400
H	5.09313300	0.40388900	-1.79452000
N	4.82059200	-1.76274400	2.28610100
H	3.86909200	-2.06582700	1.99127100
H	5.29385400	-2.54195700	2.77872200
C	4.35181100	-3.22528700	-1.42402200
H	4.44158900	-4.33060200	-1.44268900
C	4.55019600	-2.68852700	-2.85468600
H	3.83104900	-3.14901300	-3.56059000
H	5.57276900	-2.91427000	-3.21687700
H	4.40177700	-1.59230800	-2.90800200
C	1.77148300	-4.80368300	-1.98970500
N	0.87540200	-5.40932300	1.21930600
H	1.57905400	-4.80688700	0.70544200
H	1.35199500	-6.23816200	1.61688400
H	0.49779400	-4.84199400	2.00697300
C	-3.12966800	-4.31704400	1.58415300
H	-4.20221000	-4.40224700	1.84771300
C	-2.29082100	-4.51765700	2.85797900
H	-2.56081900	-3.79009800	3.64853400
H	-1.20778200	-4.35795000	2.65848600
C	-4.87158400	-1.76648600	2.07266000
H	-5.19850600	-2.79775500	2.31046700
H	-5.75321500	-1.27958900	1.61195100
C	-4.52685600	-1.02431300	3.37693700
H	-4.22906900	0.02541500	3.18359900
H	-3.68785800	-1.51531700	3.90940500
H	-5.40177400	-1.00978300	4.05650200
C	-4.90288400	0.89592500	0.43770900
C	-5.94510900	0.22471000	-0.23552900
C	-7.29572500	0.53102300	-0.05129600
H	-8.08163000	-0.00801700	-0.60317300
C	-7.63865000	1.54371800	0.86048500
H	-8.69743100	1.79428100	1.01929800
C	-6.63011700	2.22350100	1.56205400
H	-6.89691700	3.01240800	2.28054200
C	-5.28014500	1.90580200	1.34829800
H	-4.49008800	2.44522300	1.89095400
N	-5.57848800	-0.85550700	-1.18232700
H	-6.41504500	-1.34442700	-1.54881200
H	-4.95295100	-1.55708100	-0.71207100
H	-5.05176100	-0.46761500	-1.99557800
C	-4.59910100	2.14316100	-2.91496600
H	-5.62569200	2.21864200	-3.32484600
H	-4.39404100	1.07716700	-2.66131300
H	-3.88725100	2.38692200	-3.72845500
C	-4.40472500	3.05788600	-1.69305400
H	-4.47738700	4.11312600	-2.02502200
H	-5.24225400	2.92108400	-0.98397300
C	-2.98783800	-0.71227100	0.52561700
C	5.58646000	-1.29084300	1.10868800
C	6.97448100	-1.45354000	1.09163000
H	7.50062200	-1.96403900	1.91306500
C	7.69229400	-0.94742500	-0.00443400
H	8.78412300	-1.07187000	-0.04198300
C	7.01408500	-0.27962200	-1.03769700
H	7.57780400	0.12557300	-1.89083700
H	4.68661400	-0.98235200	2.97087500
C	-0.86056700	-4.77272500	-0.42258800
C	-0.20366700	-5.80022600	0.28524400
C	-1.83570200	-5.17656300	-1.35916500
C	-0.49794000	-7.15567100	0.11954300
C	-2.13831300	-6.53163500	-1.56128300
H	-2.35955100	-4.40217700	-1.93726300
C	-1.48036700	-7.52271500	-0.81552500
H	0.03167900	-7.92736100	0.69982600
H	-2.89794000	-6.81656800	-2.30388100
H	-1.72131800	-8.58557600	-0.96123100
H	1.27357200	-5.69133400	-1.55426100
H	2.80174900	-5.13715100	-2.22468100
C	1.05027700	-4.43186300	-3.29805100
H	1.55612200	-3.58950300	-3.81065300
H	1.03425300	-5.29654500	-3.99086600
H	0.00122500	-4.12613500	-3.11585100
H	-2.93861100	-5.14131200	0.87432100
H	5.18031000	-2.87857300	-0.78183400
C	0.86182300	7.72002800	-0.17058000
H	-0.68197400	7.73091300	1.35929700
H	2.38864000	7.39676400	-1.69120400
H	0.97224800	8.81383800	-0.19162200
H	-1.03428500	-0.56753700	-0.28347300
H	0.86882500	0.64574100	-0.15298400
H	-2.42339700	-5.53723600	3.26970400

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$[\alpha_2\beta_2\text{-}\mathbf{P}]^{6+}$

C	-0.89135700	2.93026300	-0.43953800
C	3.94642500	-1.00485800	1.12461100
C	4.07307000	0.39874100	1.11185000
C	2.93622800	0.93119500	0.37892400
C	2.73063200	-1.34966700	0.41537400
C	2.24876700	-2.64661100	0.05104400
C	-3.94365600	1.09281800	1.11061900
C	-4.06908600	-0.31077500	1.11353500
C	-2.93122600	-0.86252300	0.40552200
C	-2.67844600	-2.22445700	0.04753600
C	-1.39622300	-2.75316100	-0.27486200
C	-1.05451700	-3.98267600	-0.97520200
C	0.35151800	-4.10072600	-0.97731200
C	0.89586900	-2.94764500	-0.27499400
C	1.04696900	3.93071500	-1.18188200
C	1.38264700	2.73449000	-0.43845600
C	2.66419900	2.26772800	-0.02126400
C	-2.23382700	2.68967700	-0.02228400
N	0.18262200	2.10383600	-0.09439300
H	0.12910200	1.49608600	0.72968500
C	-0.35568100	4.05167800	-1.18227300
C	1.99559400	4.79692500	-1.96014800
H	2.84276300	5.10866000	-1.32196500
H	1.48197900	5.73685200	-2.23957900
N	-2.10855800	0.20742800	0.05033900
H	-1.44214600	0.13767600	-0.72418700
N	-0.18945000	-2.13429000	0.07009800
H	-0.13002600	-1.41430100	0.79444900
N	2.10816500	-0.15377100	0.05786000
H	1.43817700	-0.10987500	-0.71525700
C	2.51086000	4.11373600	-3.24035100
H	1.70720700	3.96555400	-3.98767900
H	2.90081000	3.09044300	-3.03275900
H	3.30208000	4.71961400	-3.72268200
C	3.72889600	3.30928000	0.12329600
C	4.92253600	3.36255100	-0.63352900
C	5.89118600	4.35032100	-0.43868100
C	4.51291700	5.30609500	1.31983300
C	3.54455500	4.31509600	1.09863800
H	2.62570000	4.30659500	1.70225000
N	5.15291400	2.37955000	-1.72129300
H	4.52327100	2.57003200	-2.53427300
H	4.98515800	1.40063000	-1.41679800
H	6.12916900	2.43075100	-2.07290900
C	5.09097200	1.16738400	1.90847000
H	5.49284500	2.02291200	1.33623400
H	5.95811000	0.50727200	2.10399800
C	4.52694100	1.67643600	3.25040500
H	3.66453100	2.35605700	3.10504800
H	5.30994700	2.23420800	3.79940600
H	4.19477300	0.83996100	3.89543100
C	4.81545700	-1.94515300	1.91090900
H	5.12003700	-2.80373700	1.28433300
H	5.75857800	-1.43127400	2.17761500
C	4.13992900	-2.43438300	3.20591400
H	4.01627000	-1.61916200	3.94513300
H	4.73985500	-3.22796800	3.69250500
H	3.10463000	-2.80348900	3.01926800
C	3.31298100	-3.68509800	-0.14531600
C	4.25336500	-3.47372800	-1.17846700
H	4.16978800	-2.56293900	-1.78727200
N	2.56497800	-5.12393400	1.76498700
H	1.56487600	-5.04963600	1.49968100
H	2.71328700	-6.07531000	2.15570900
C	1.11241600	-5.12135200	-1.77905500
H	0.45591900	-5.99538500	-1.95320600
C	1.59123400	-4.57292600	-3.13902800
H	0.74023900	-4.25803900	-3.77304000
H	2.14552700	-5.35989400	-3.68584600
H	2.26488600	-3.70218000	-3.02115700
C	-1.97801000	-4.86548200	-1.76906200
N	-3.44509600	-4.57464400	1.80462300
H	-2.44497900	-4.70113000	1.56106700
H	-3.77718500	-5.46642500	2.22157500
H	-3.50409500	-3.84317400	2.55183100
C	-5.09520900	-1.08559600	1.89043300
H	-5.94436100	-0.42011000	2.13718100
C	-4.53045800	-1.66953800	3.19979000
H	-4.30699100	-0.88193700	3.94542400
H	-3.55457300	-2.18447700	3.03805700
C	-4.80050300	2.02688900	1.91990800
H	-5.76270400	1.52518500	2.13922800
H	-5.06323200	2.93492600	1.34730700
C	-4.12935300	2.44169100	3.24510500
H	-3.16092700	2.95151900	3.07477300
H	-3.94176000	1.56781300	3.89891600
H	-4.78838800	3.13913500	3.79740400
C	-3.10205400	3.89919700	0.12557600
C	-4.26944500	4.16021100	-0.62905100
C	-5.05107400	5.30097500	-0.43081300
H	-5.94875800	5.48496600	-1.04179900
C	-4.67886500	6.22442400	0.56237900
H	-5.29496000	7.12042600	0.72672700
C	-3.52387800	6.00015900	1.32601900
H	-3.22398500	6.72107800	2.10053400
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H	-5.61612000	3.45940200	-2.07651400
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C	4.47703600	-5.79869900	0.36767900
H	4.57500600	-6.70437400	0.98668100
C	5.38141100	-5.56869400	-0.68409600
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H	5.97519200	-4.21132600	-2.27489000
H	2.74350800	-4.44029100	2.53749900
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C	-4.77393000	-2.72656800	-1.20570100
C	-5.41336900	-4.94882800	0.37464600
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H	6.80553400	4.37497400	-1.05239500
H	6.45088500	6.10053600	0.71655400
H	-5.25174500	-2.36623500	3.66975800

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$\alpha_2\beta_2\text{-}\mathbf{P}[H_2SO_4 \cdot H_2O]_2]^{4+}$

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C	-1.63251808	-3.21515661	-1.23237725
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H	4.16766197	4.74227771	-3.09719365
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C	7.02979888	2.48363038	-0.72996292
H	7.79293305	2.14932616	-1.45026201
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C	6.35176148	3.94064828	1.09673799
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N	5.48776545	0.77528129	-1.63169654
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H	5.45036939	0.80036551	4.05032748
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H	1.09459774	-3.71863790	-4.73970998
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H	-5.10468467	0.14880276	3.43113285
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H	-3.20092512	0.29230937	5.10178769
H	-2.53206297	-1.24920463	4.47842858
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H	-2.38250488	7.21694471	2.57861040
C	-1.78545244	5.26140725	1.85427361
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N	-3.44483285	3.23576975	-0.85807406
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C	-6.47645792	-3.64634994	0.92125459
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H	-7.28211425	-1.71590762	0.32630052
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H	-2.62989587	-4.21192541	-2.84645677
C	-3.25820773	-2.12467052	-2.80566838
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H	-3.54211251	-1.34066235	-2.07442093
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S	0.55433729	-0.87053973	3.37431097
O	1.71926909	-1.20852454	4.20740659
O	-0.50723048	-1.95145771	3.26532936
O	0.85768858	-0.34994342	1.97985224
O	-0.39088717	-4.31962910	2.40123476
H	-0.29516664	-4.81708706	3.24215141
H	-0.31183851	-3.32973591	2.72463185
H	0.90021050	-5.04205168	-2.61951618
S	-0.27906186	1.13535521	-2.46787809

Experimental NMR Data

$\alpha_2\beta_2$-P·10CSA(S)

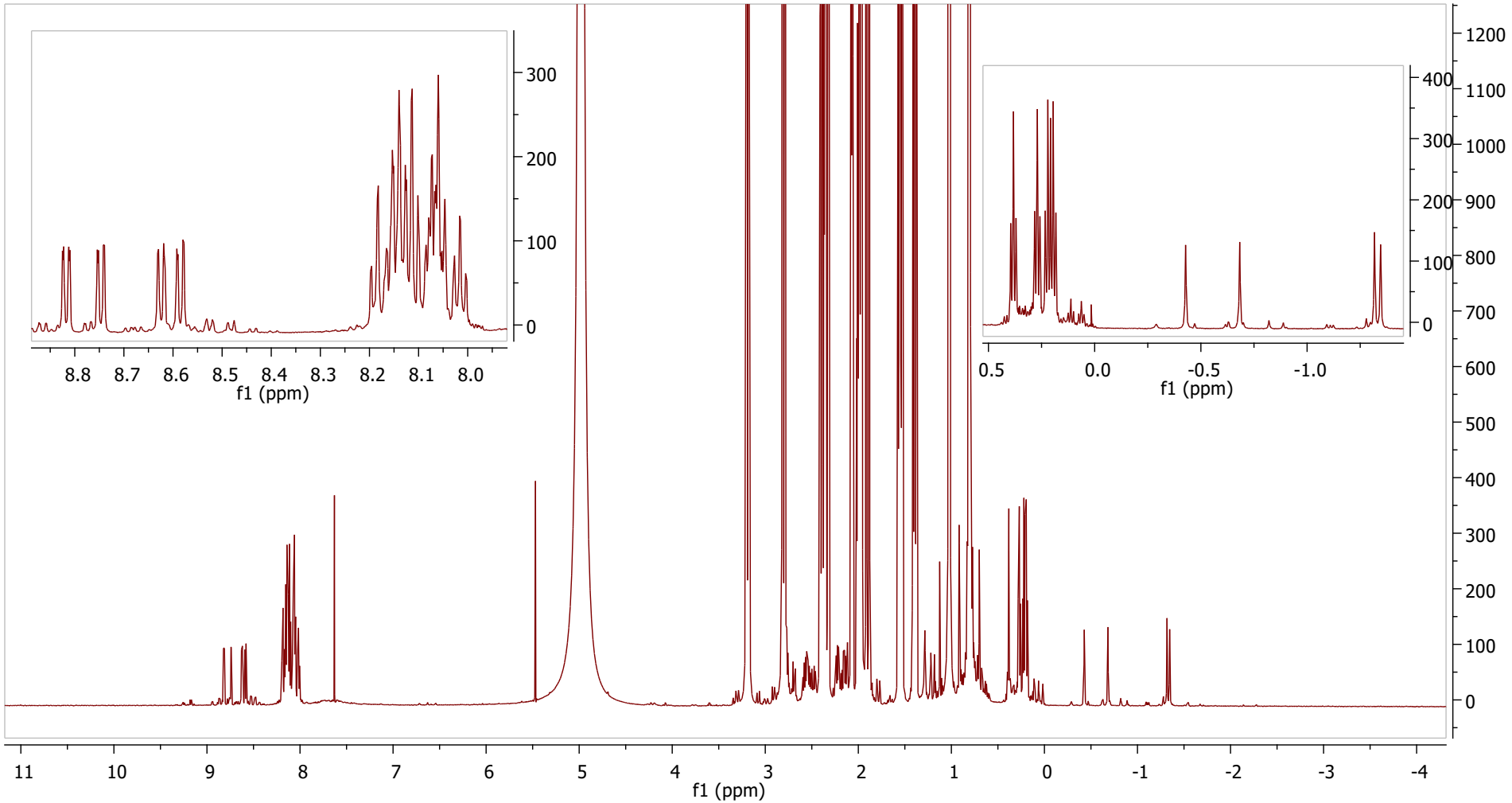

Figure S16. ^{1}H NMR spectrum of $\alpha_2\beta_2$-**P·10CSA(S)** with the expansion of areas of interest (600 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

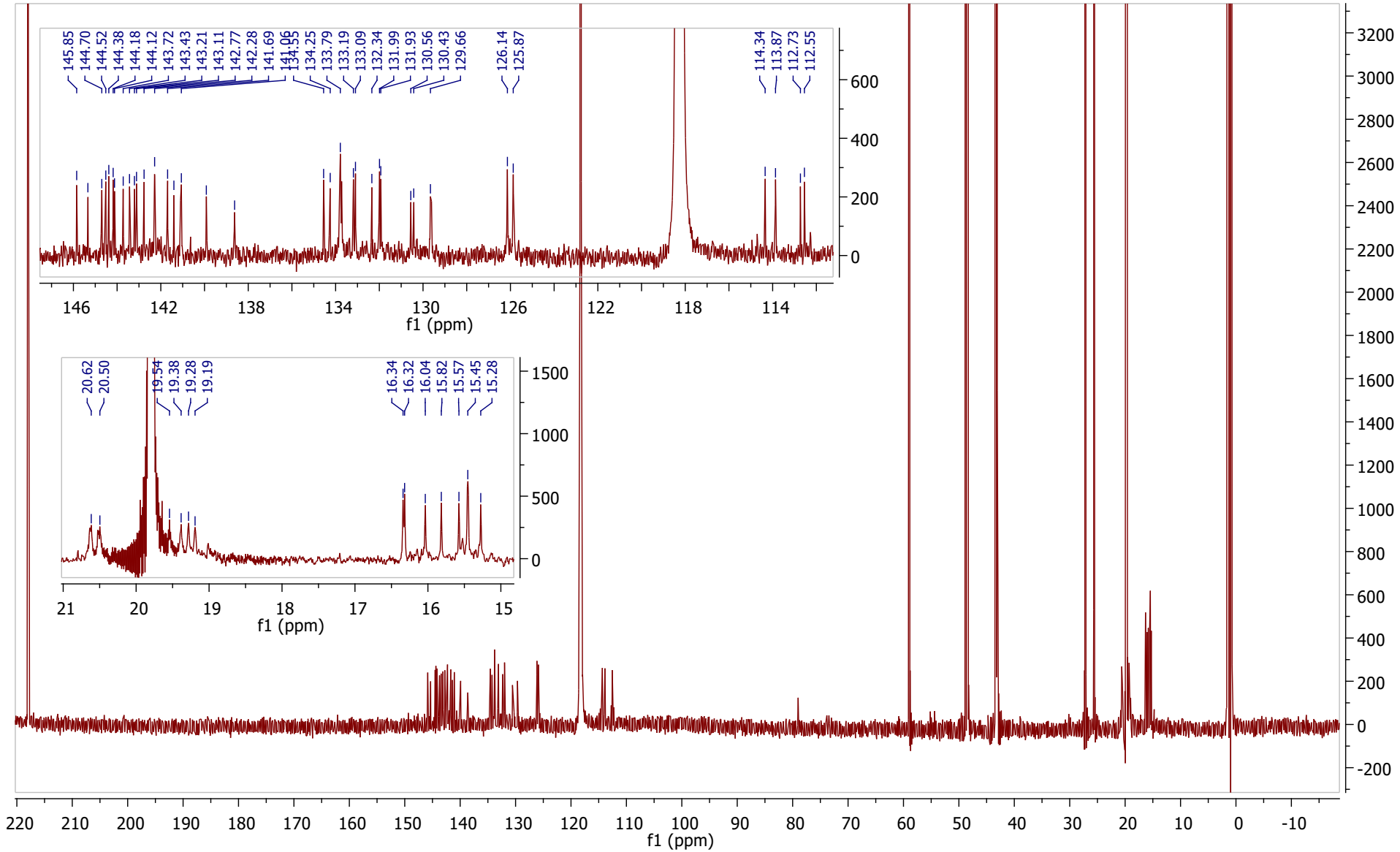

Figure S17. ^{13}C NMR spectrum of $\alpha_2\beta_2$-**P·10CSA(S)** with expansion of areas of interest (151 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

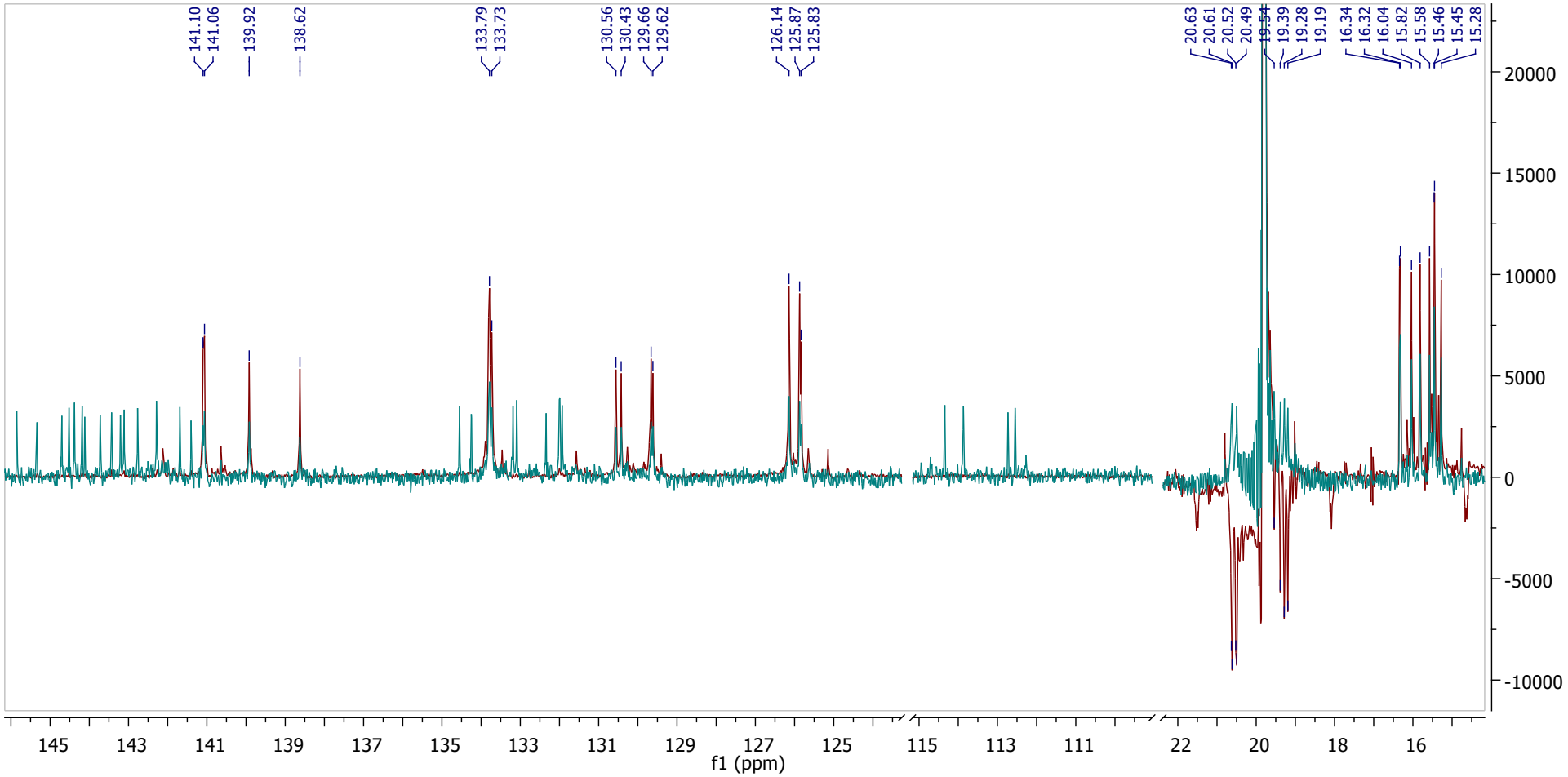

Figure S18. ^{13}C NMR (magenta) and DEPT-135 (red) overlay spectra of $\alpha_2\beta_2$-**P·10CSA(S)** (151 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

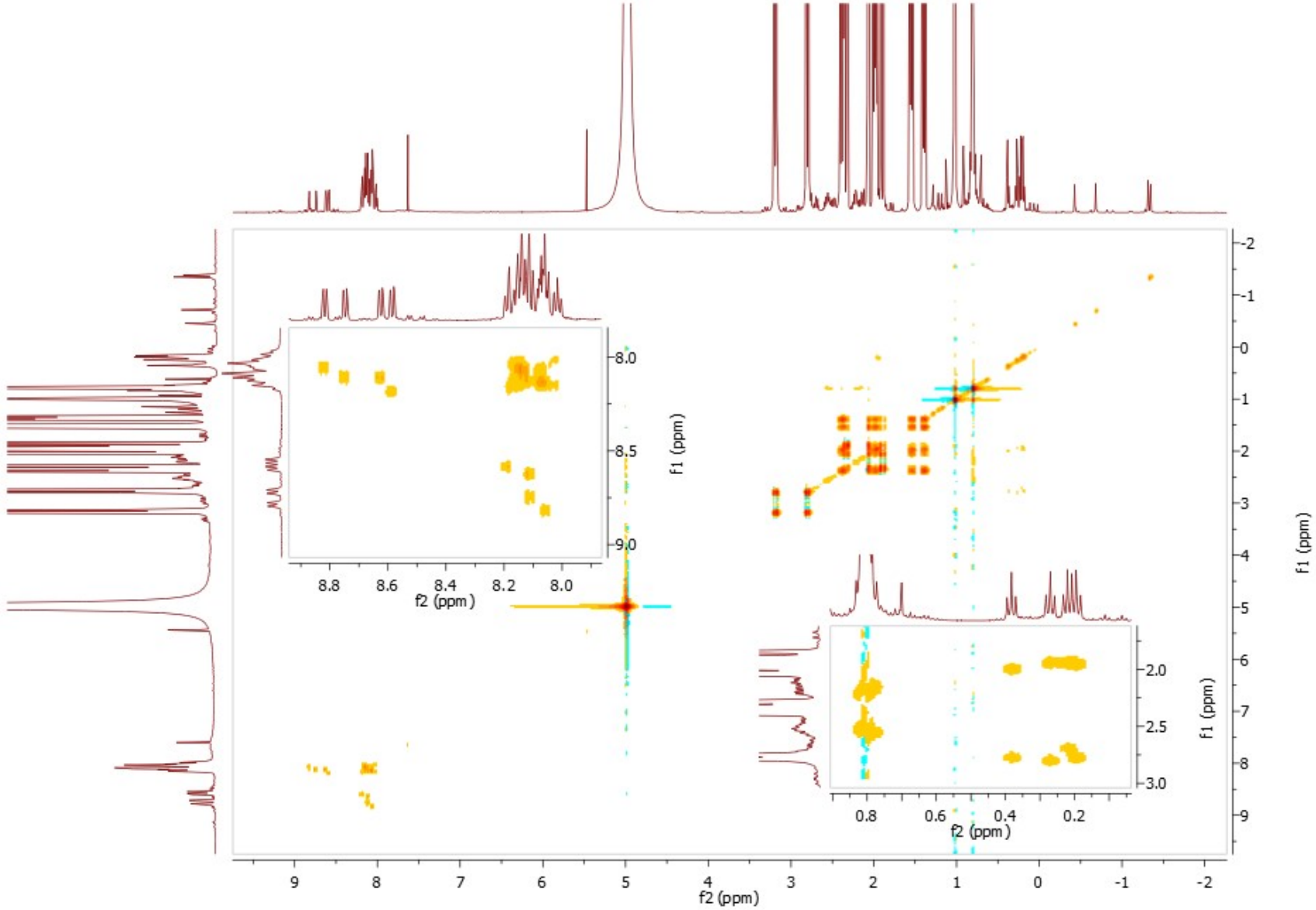

Figure S19. ^{1}H–^{1}H TOCSY spectrum of α_2,β_2-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

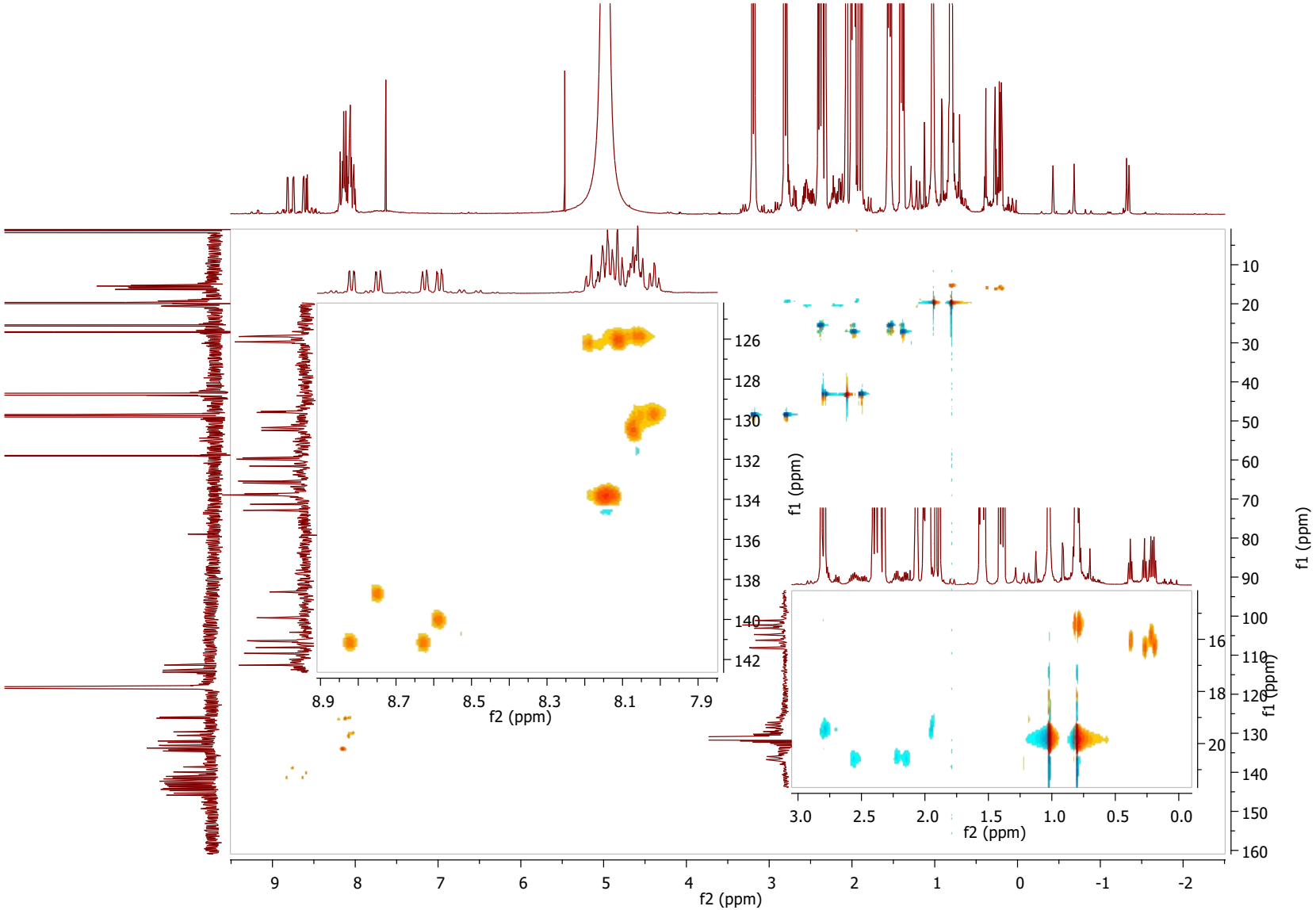

Figure S20. ^{1}H–^{13}C HSQC spectrum of α_2,β_2-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

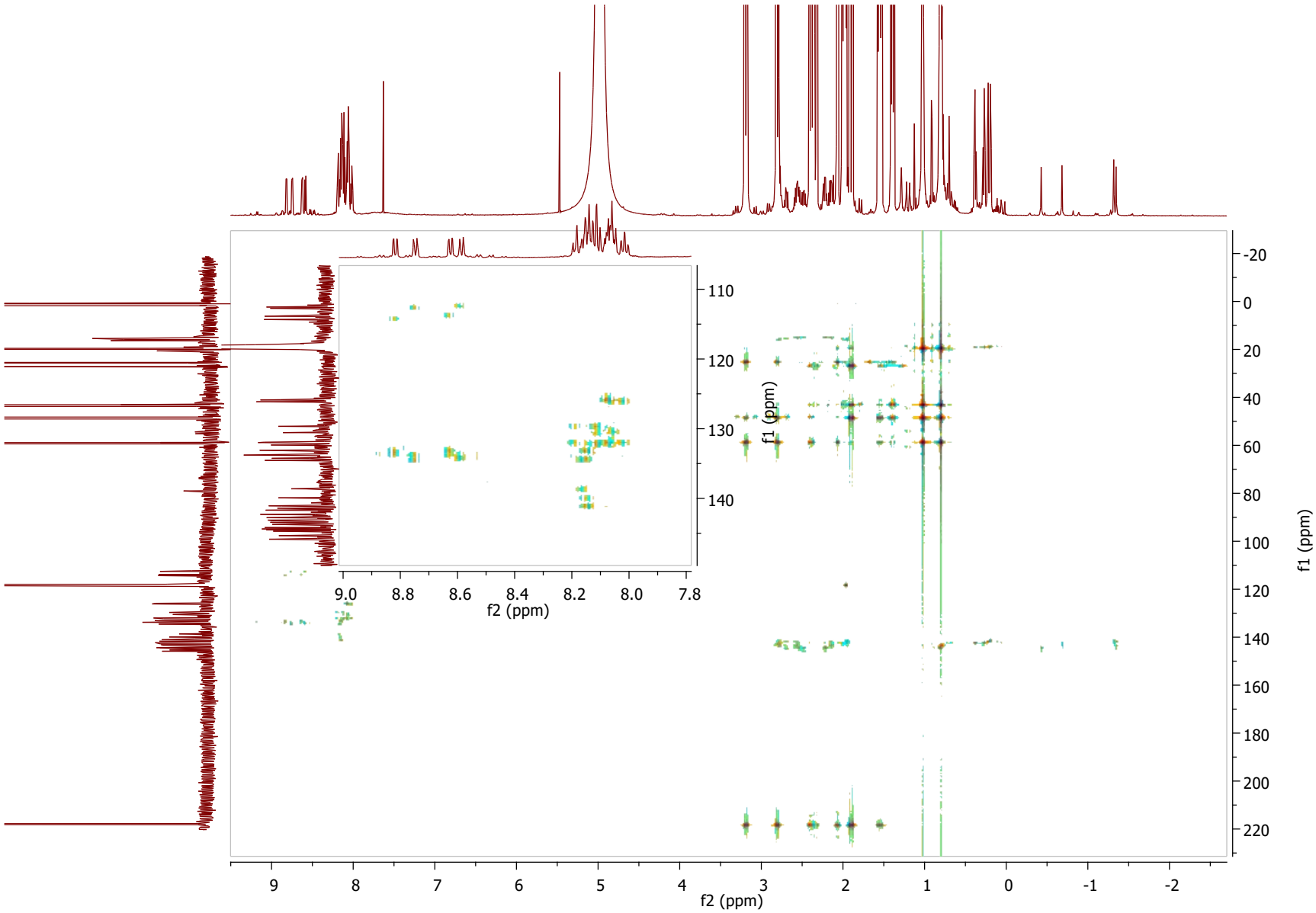
110
120
130
140
9.0
8.8
8.6
8.4
8.2
8.0
7.8
f2 (ppm)
f1 (ppm)
-20
0
20
40
60
80
100
120
140
160
180
200
220
f1 (ppm)
9
8
7
6
5
4
3
2
1
0
-1
-2
f2 (ppm)

Figure S21. $^{1}H–^{13}C$ HMBC spectrum of α_2,β_2-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

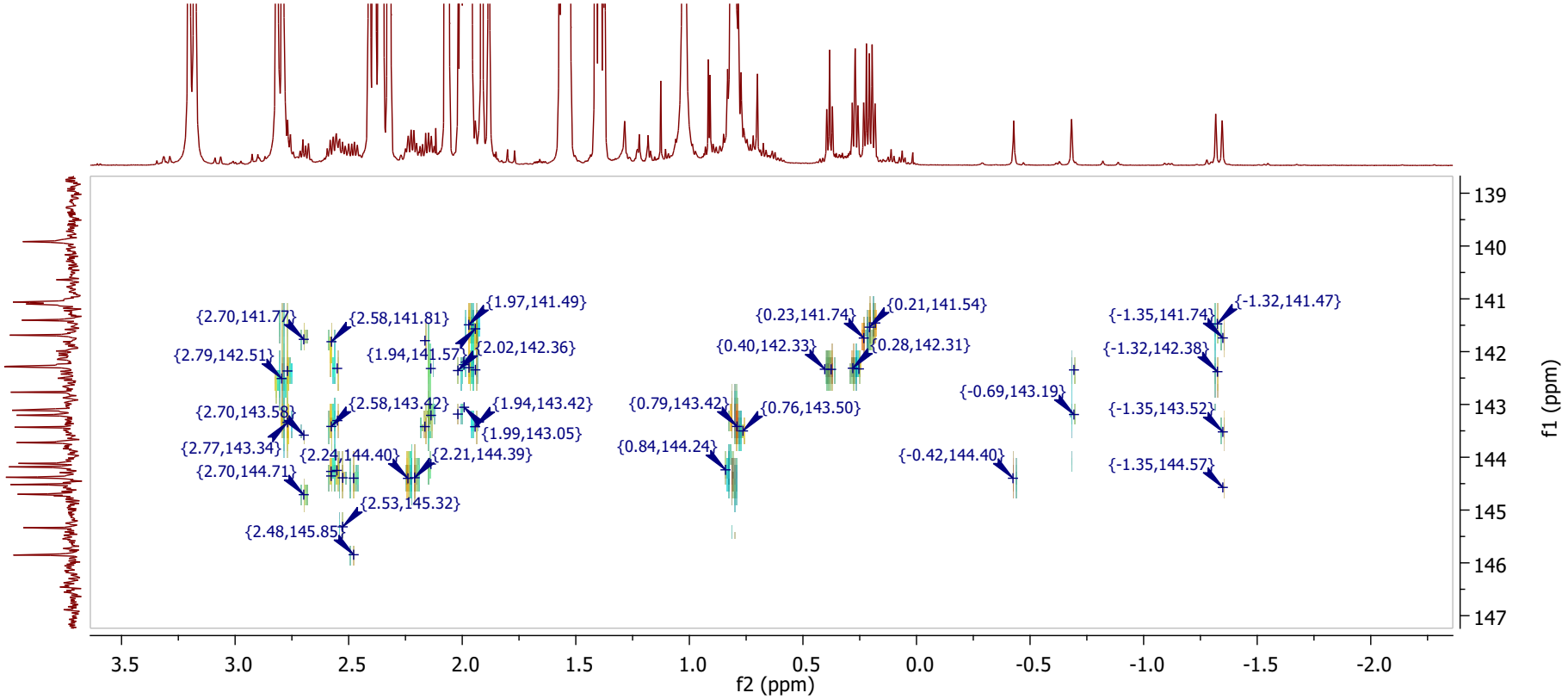

Figure S22. $^{1}H–^{13}C$ HMBC spectrum, expansion of area of interest of α_2,β_2-**P·10CSA(S)** (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

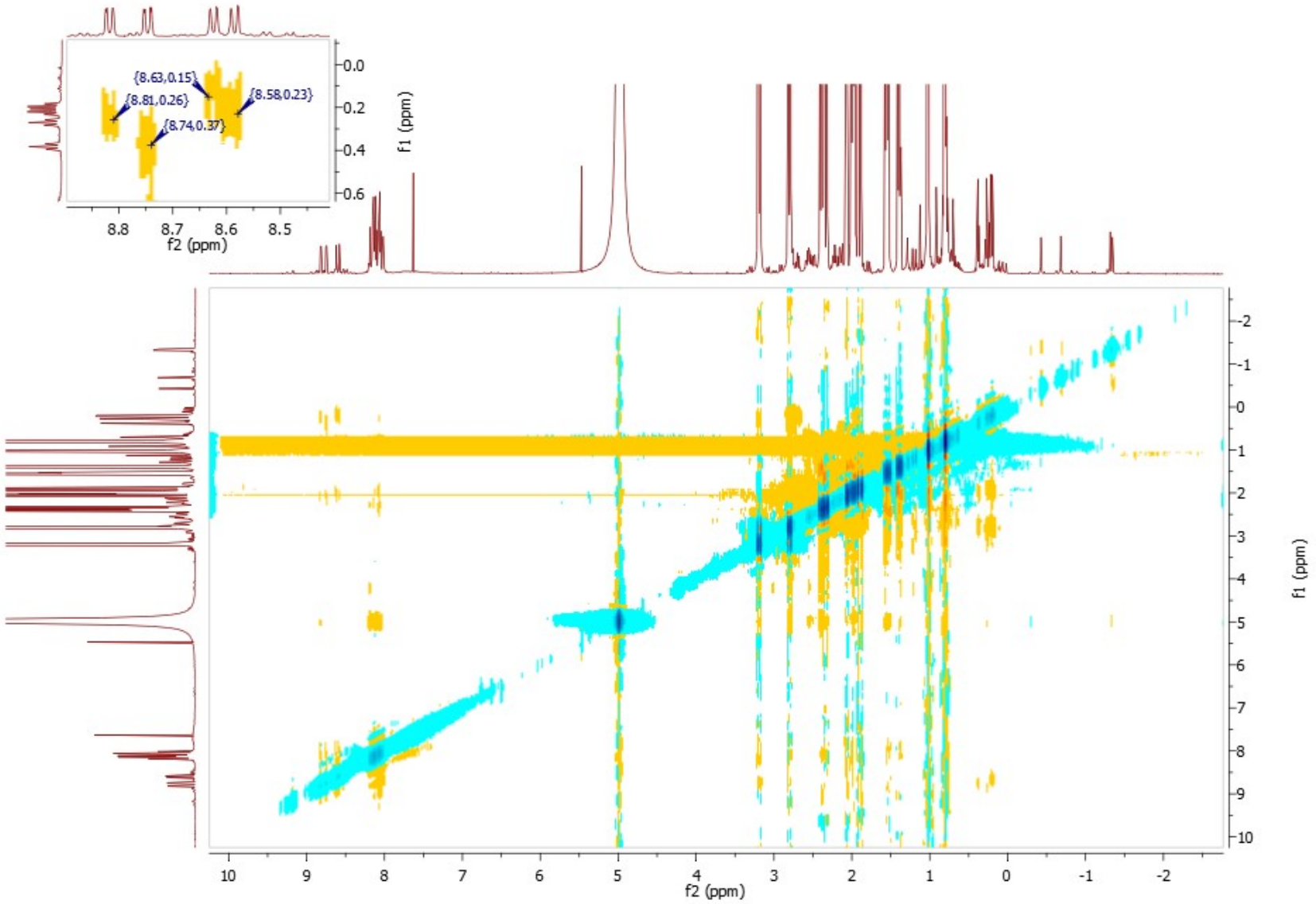

Figure S23. ^{1}H–^{1}H ROESY spectrum of $\alpha_2\beta_2$-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

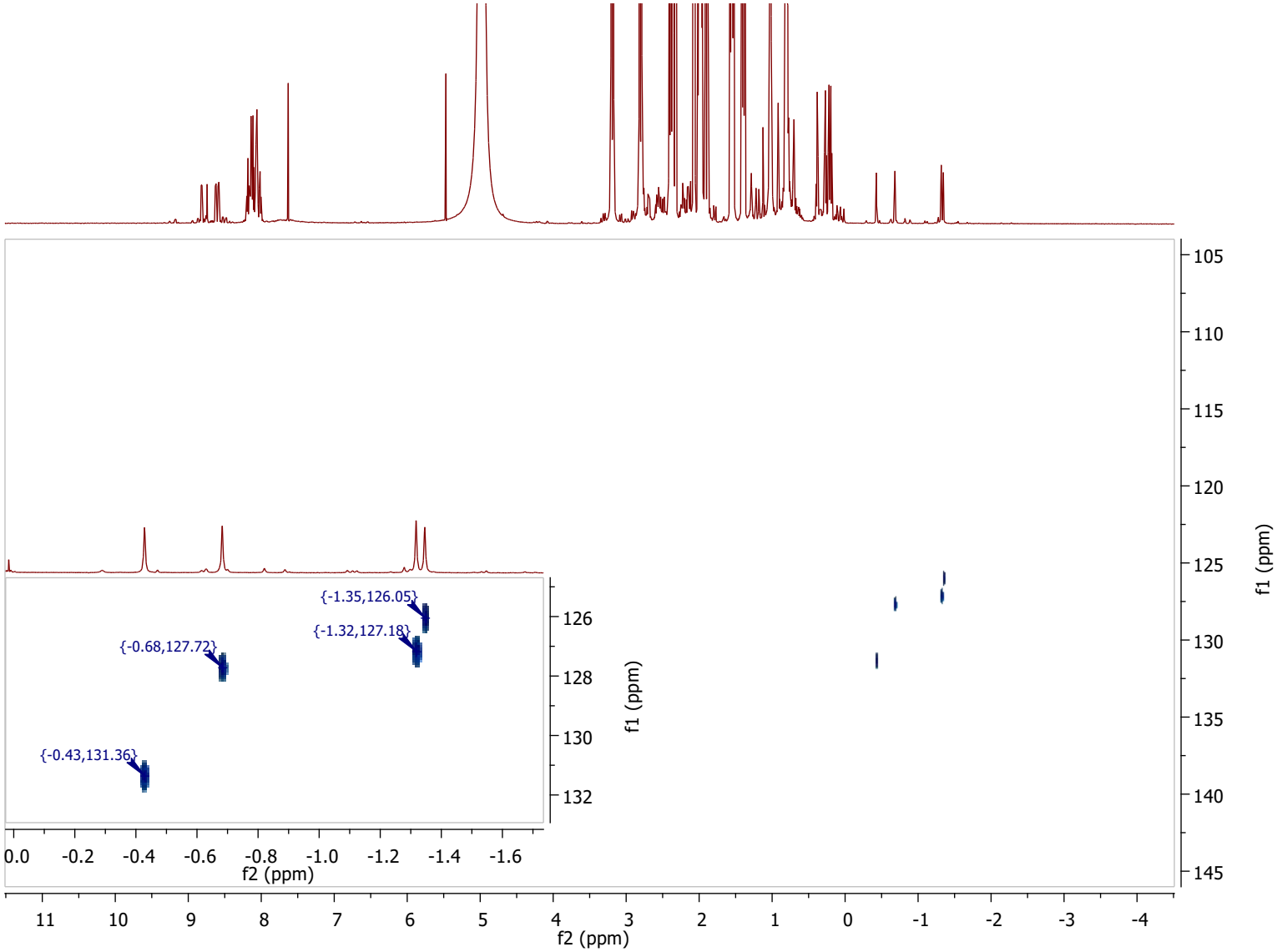

Figure S24. $^{1}H–^{15}N$ HSQC spectrum of α_2,β_2-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

$\alpha_2\beta_2$-P-10CSA(SR)

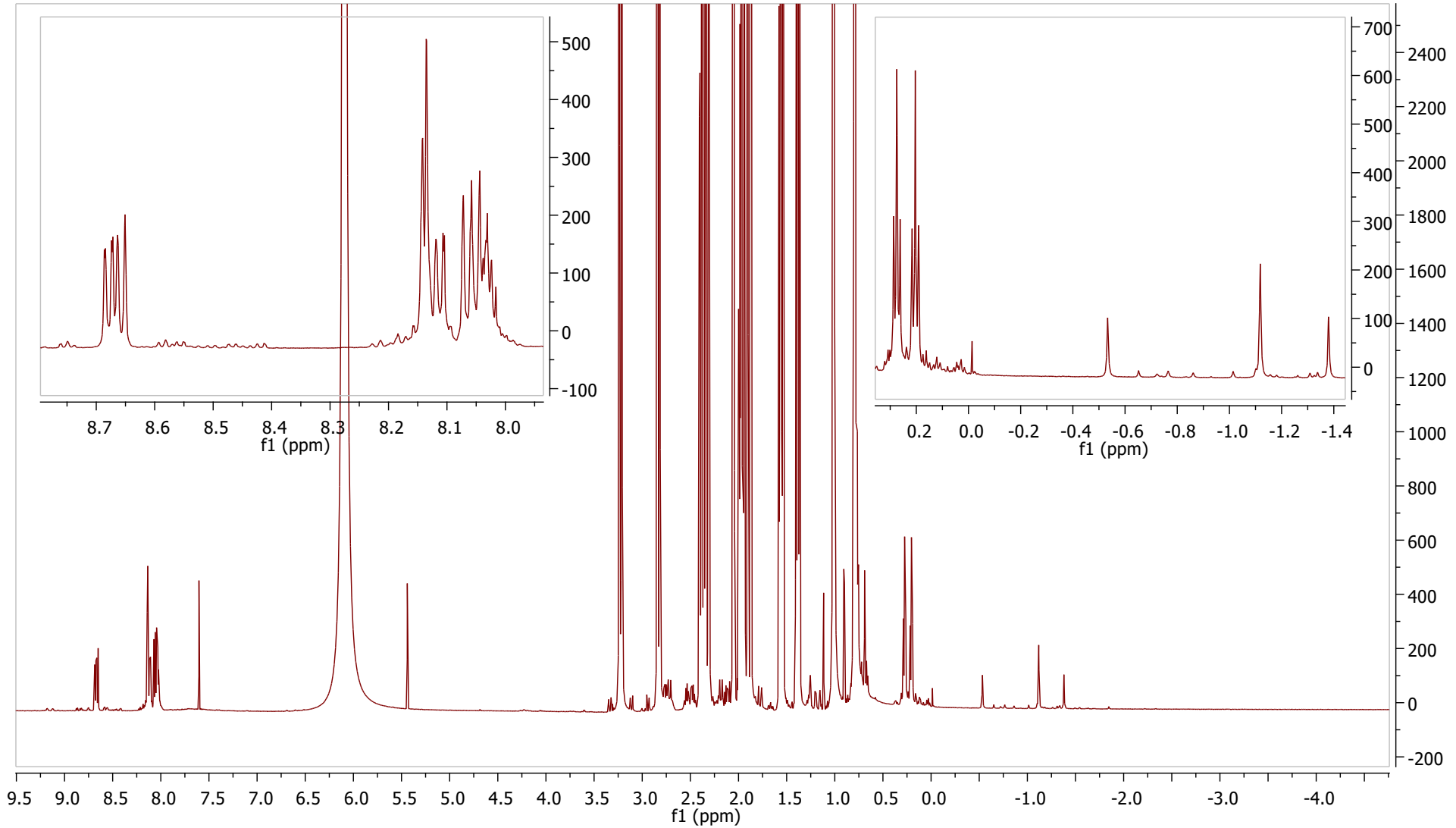

Figure S25. ^{1}H NMR spectrum of α_2,β_2-**P·10CSA(SR)** with expansion of areas of interest (600 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

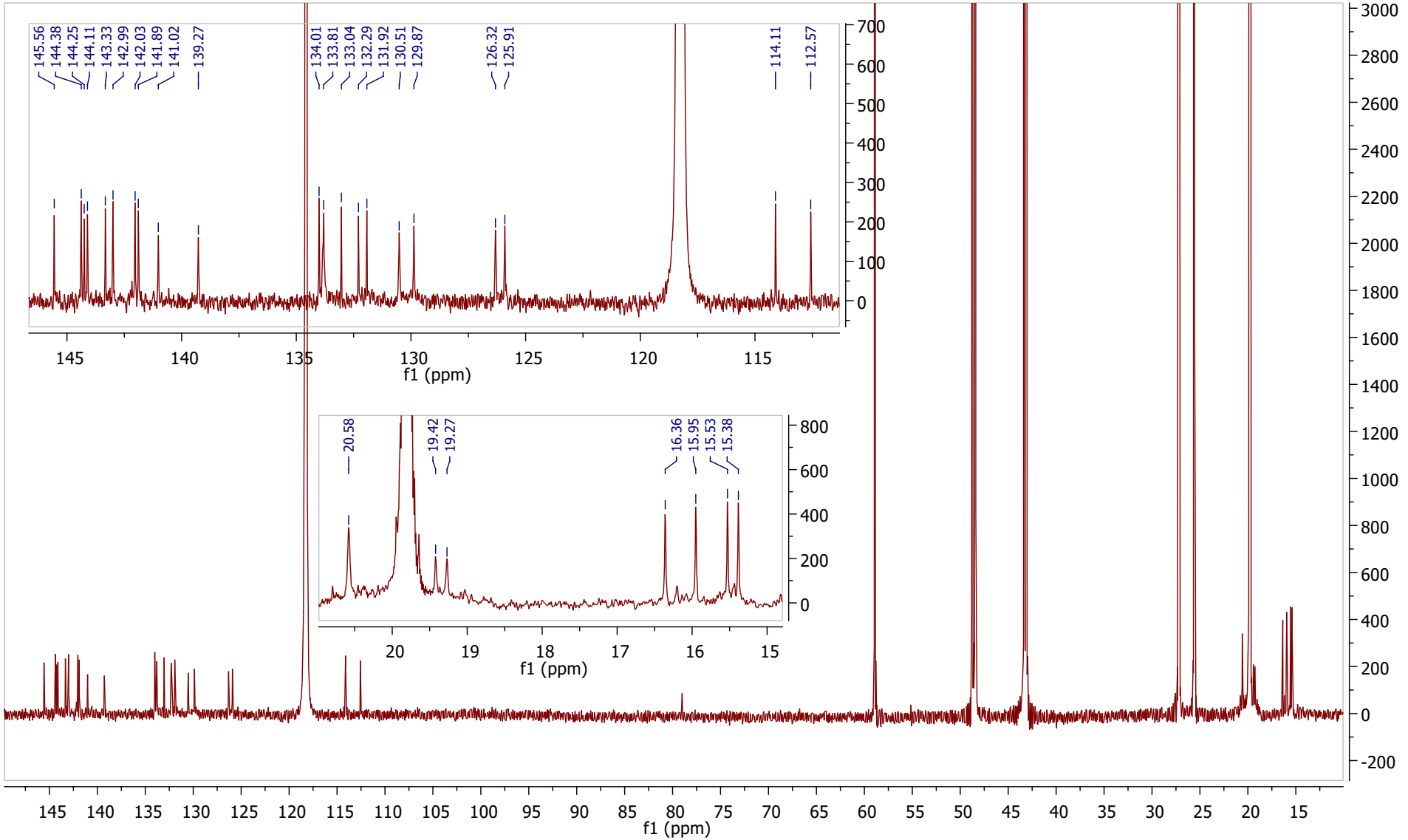

Figure S26. ^{13}C NMR spectrum of α_2,β_2-**P·10CSA(SR)** with expansion of areas of interest (151 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

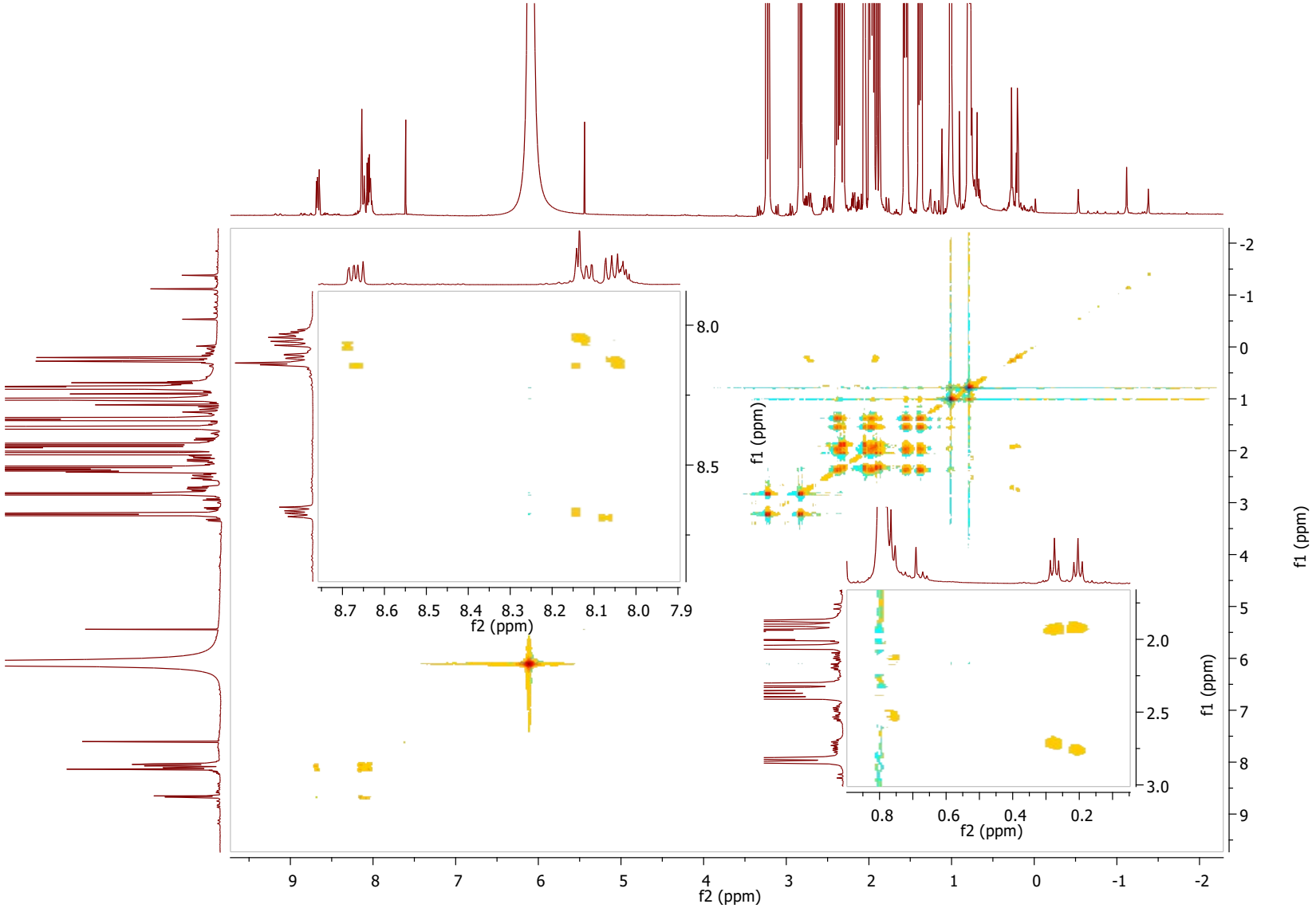
f2 (ppm)
f1 (ppm)
8.7 8.6 8.5 8.4 8.3 8.2 8.1 8.0 7.9
8.0
8.5
0.8 0.6 0.4 0.2
2.0
2.5
3.0
9 8 7 6 5 4 3 2 1 0 -1 -2
-2 -1 0 1 2 3 4 5 6 7 8 9

Figure S27. $^{1}H–^{1}H$ TOCSY spectrum of $\alpha_{2,}\beta_{2}$-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

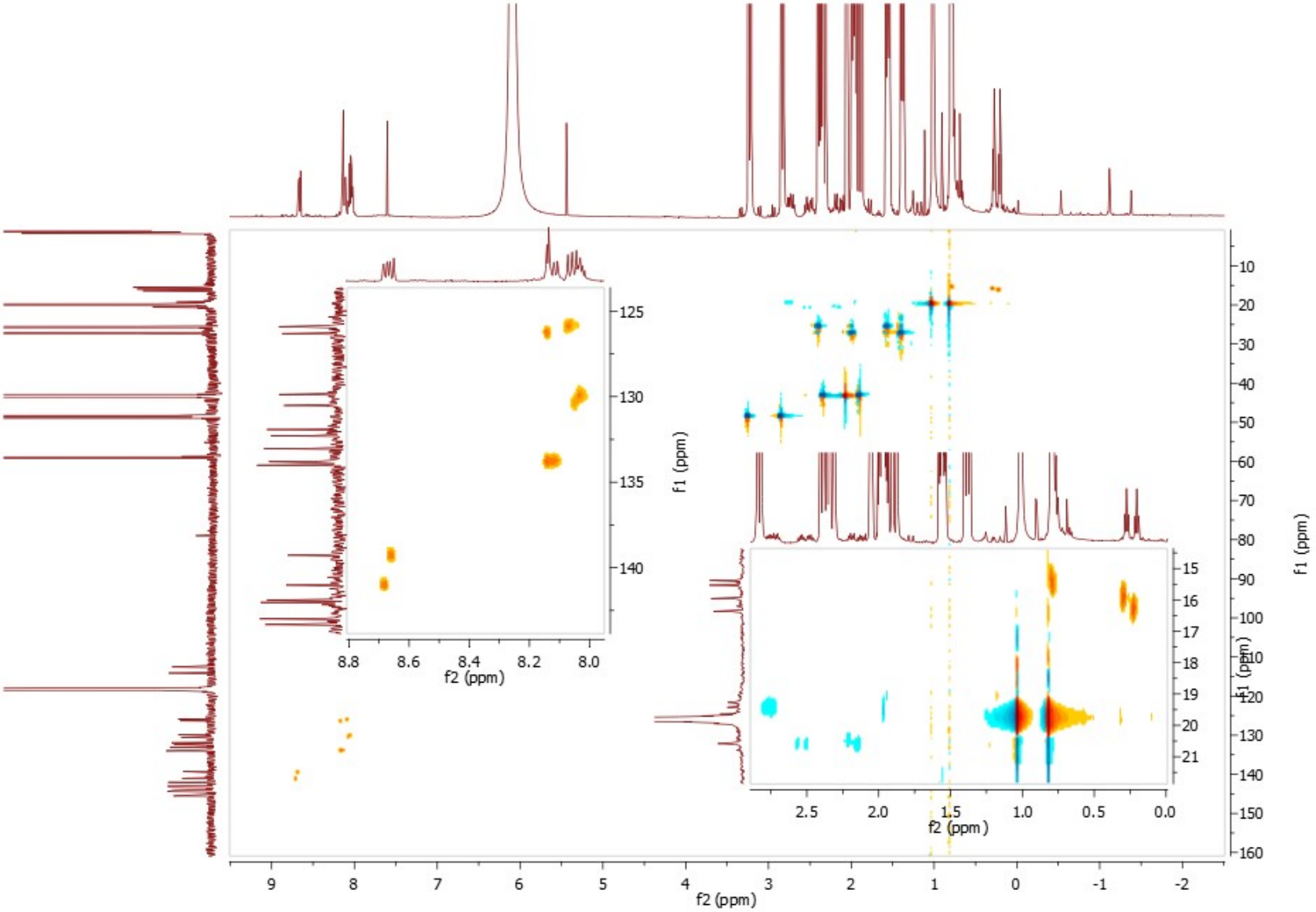

Figure S28. $^{1}H–^{13}C$ HSQC spectrum of α_2,β_2-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

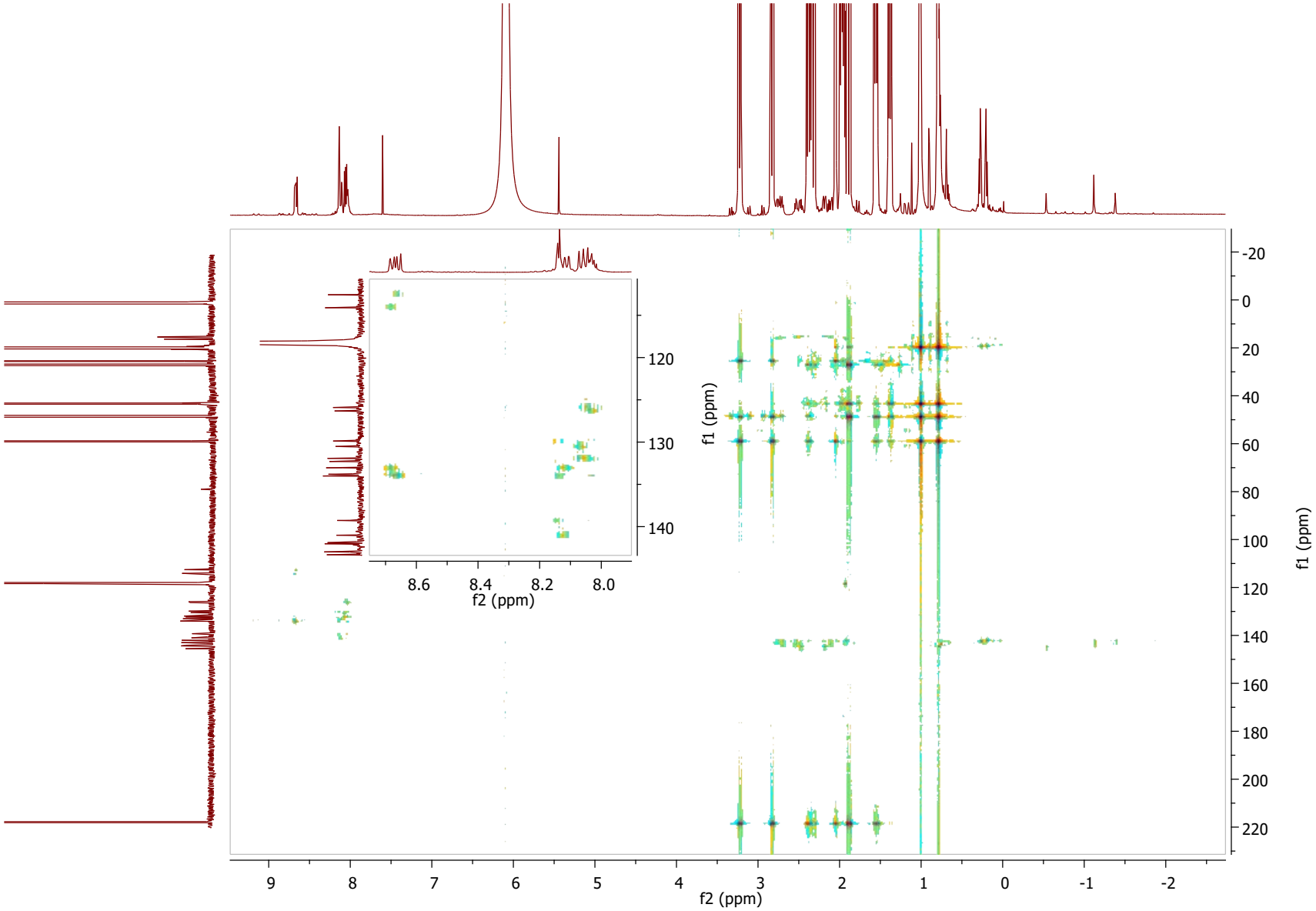

Figure S29. $^{1}H–^{13}C$ HMBC spectrum of α_2,β_2-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

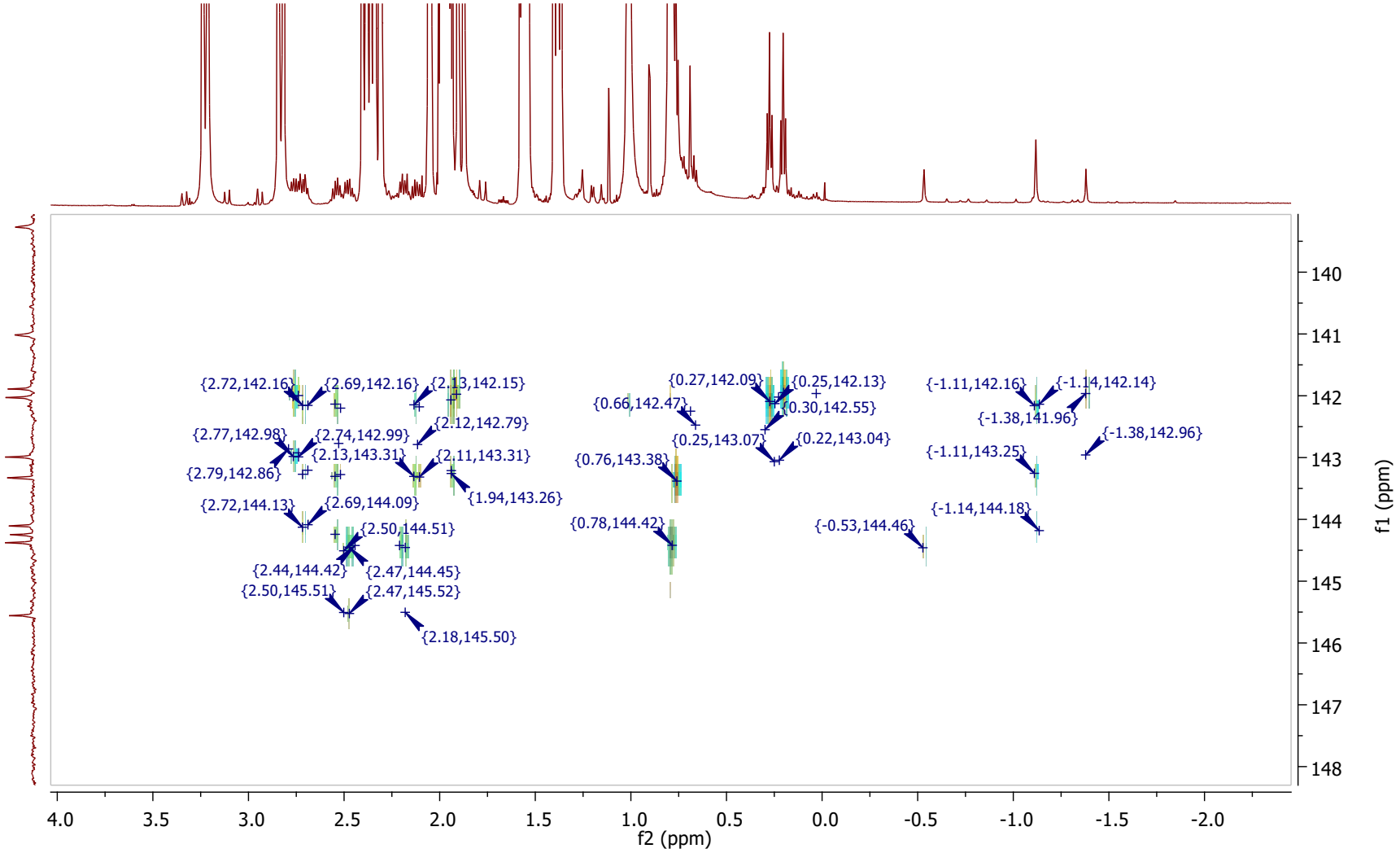

Figure S30. $^{1}H–^{13}C$ HMBC spectrum, expansion of area of interest of α_2,β_2-**P·10CSA(SR)** (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

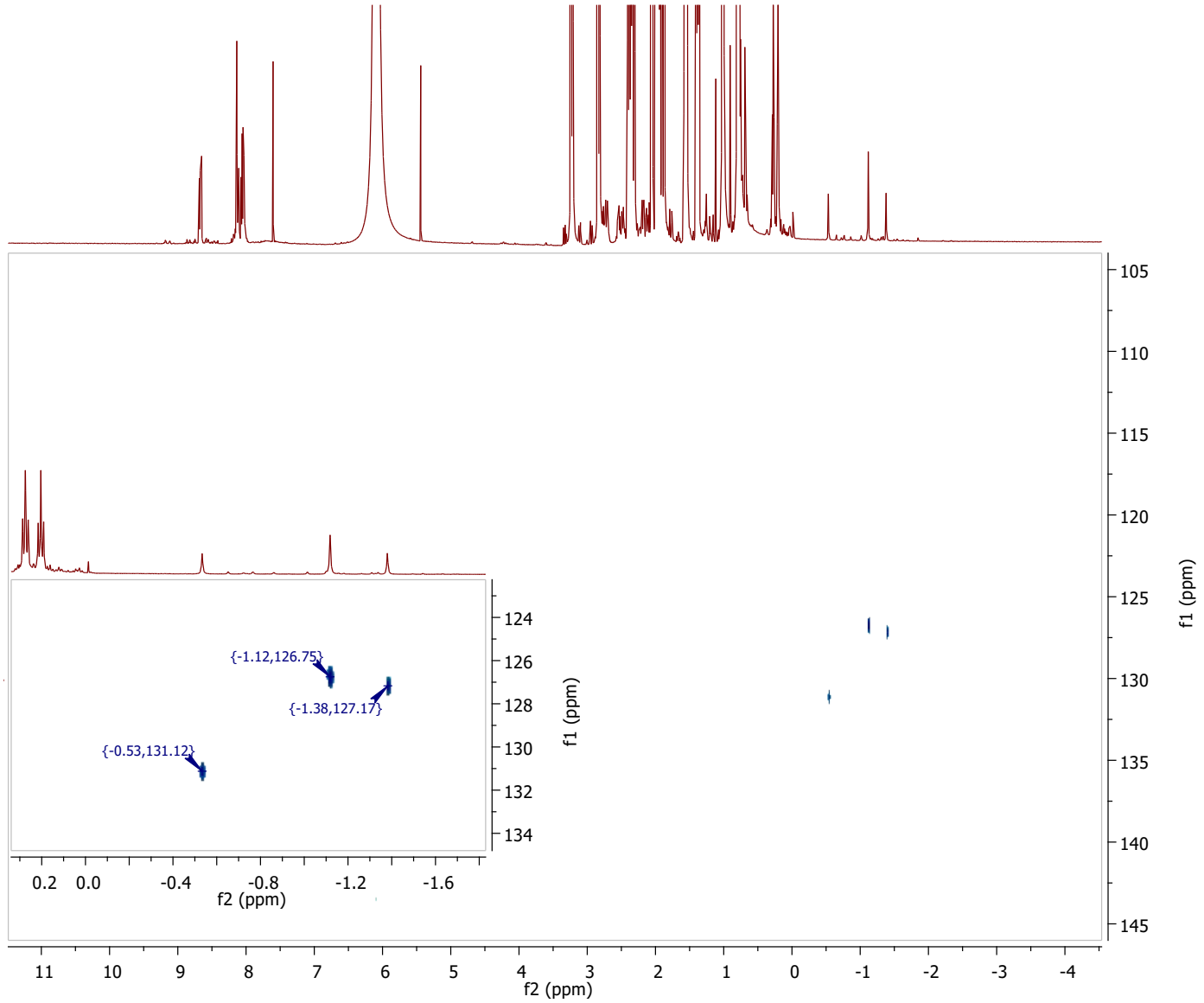

Figure S31. $^{1}H–^{15}N$ HSQC spectrum of α_2,β_2-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C)

α_4-P-10CSA(S)

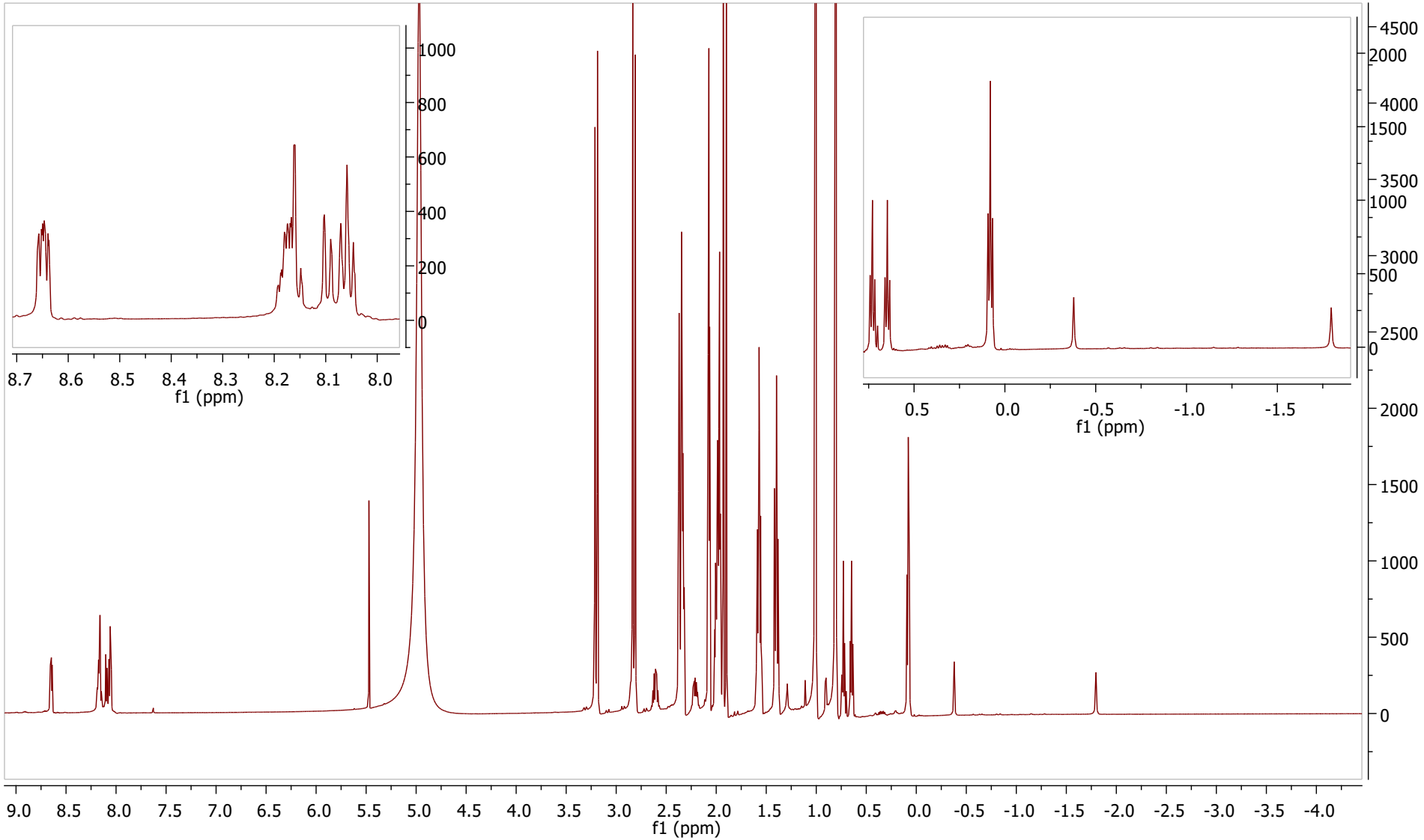

Figure S32. ^{1}H NMR spectrum of **α_4-P·10CSA(S)** with the expansion of areas of interest (600 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

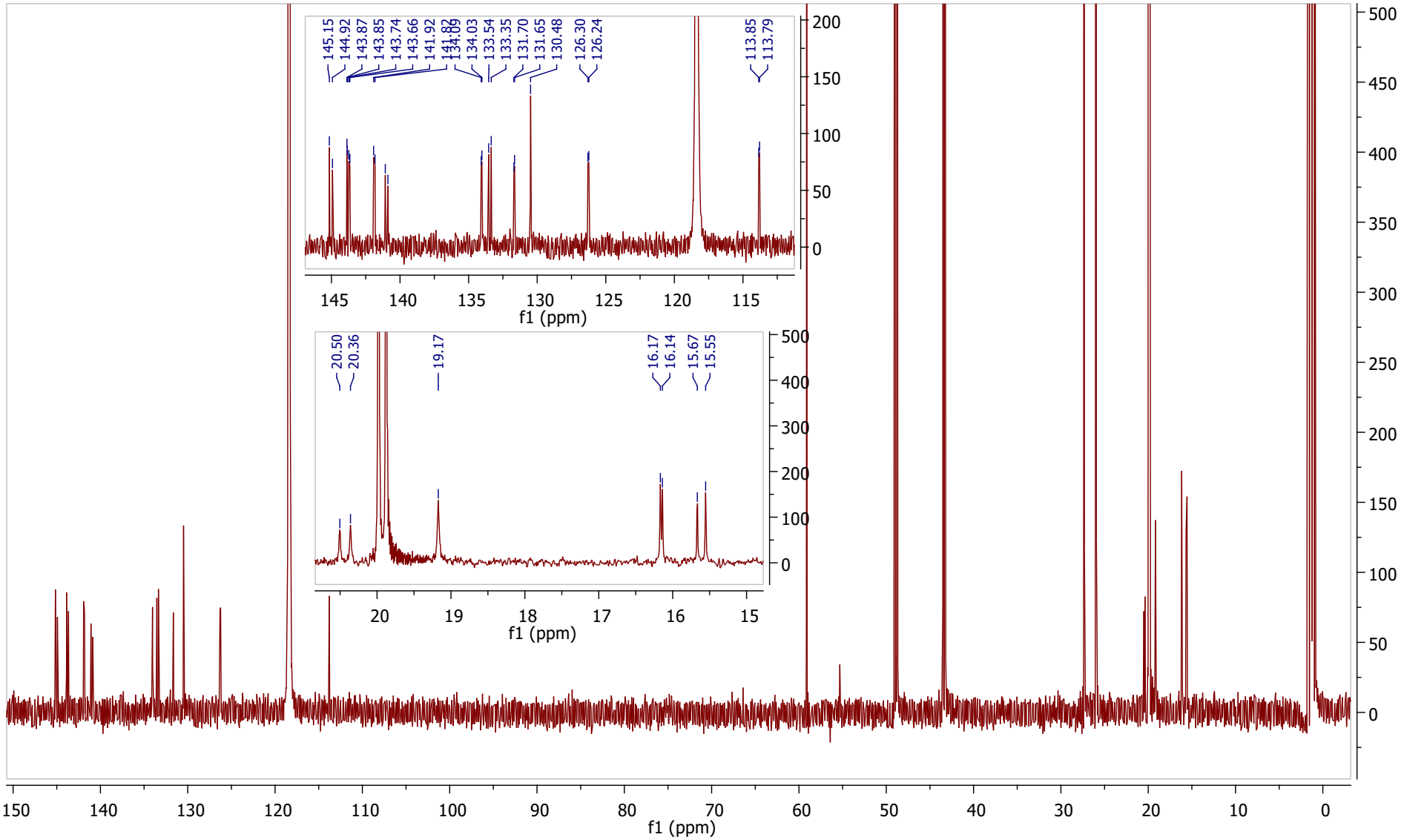

Figure S33. ^{13}C NMR spectrum of α_4-**P·10CSA(S)** with expansion of areas of interest (151 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

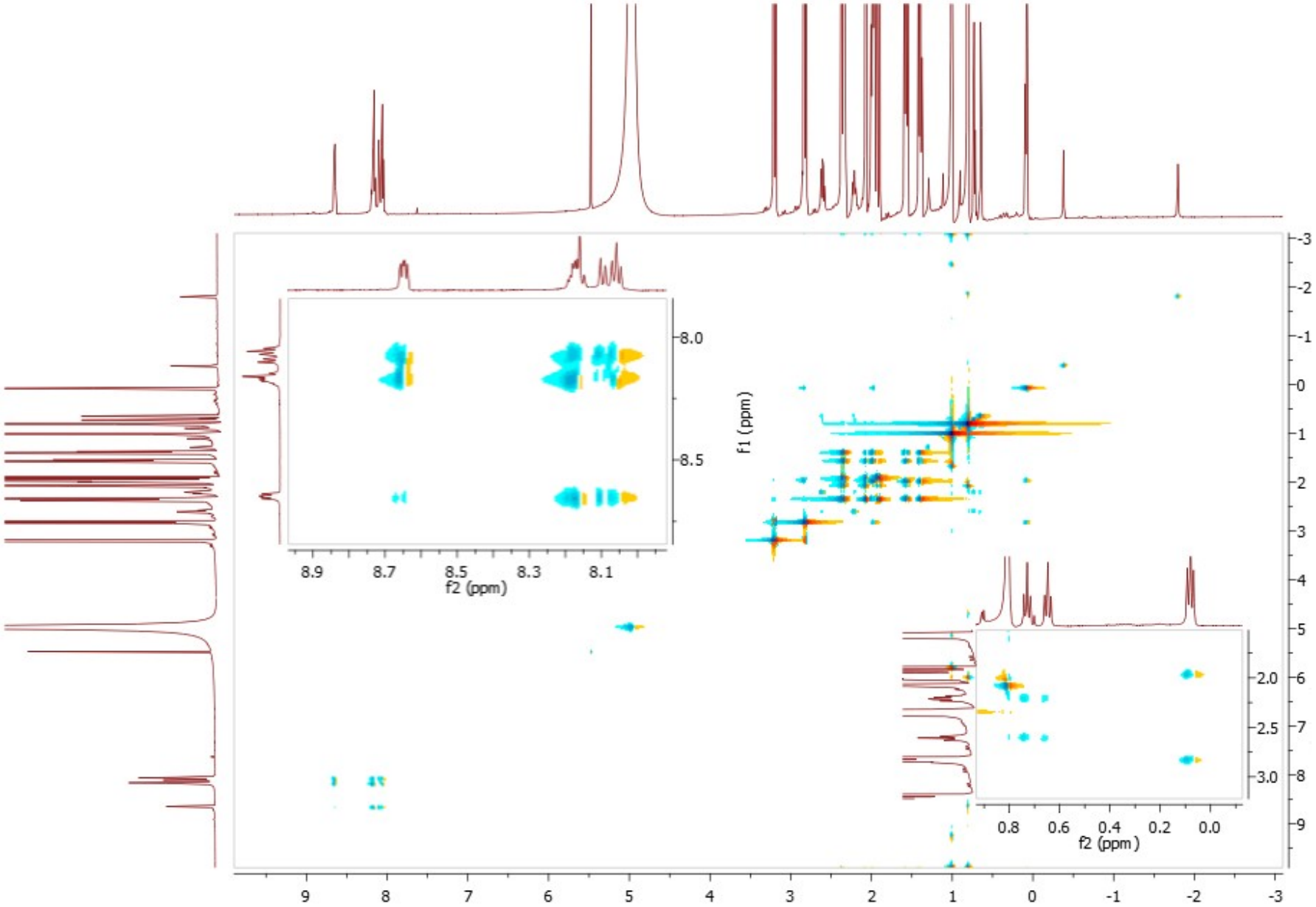

Figure S34. $^{1}H–^{1}H$ TOCSY spectrum of α_4-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

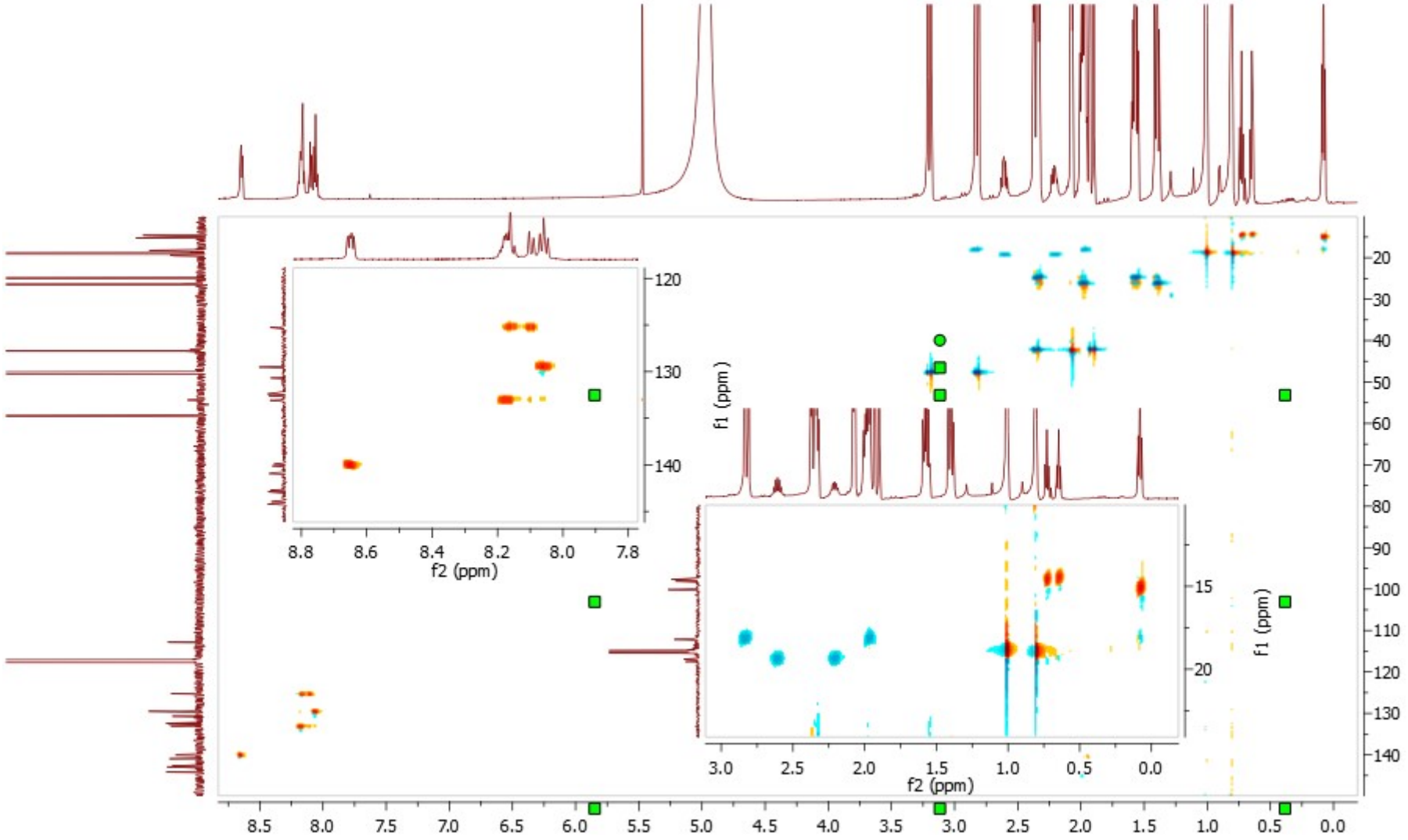

Figure S35. $^{1}H–^{13}C$ HSQC spectrum of α_4-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

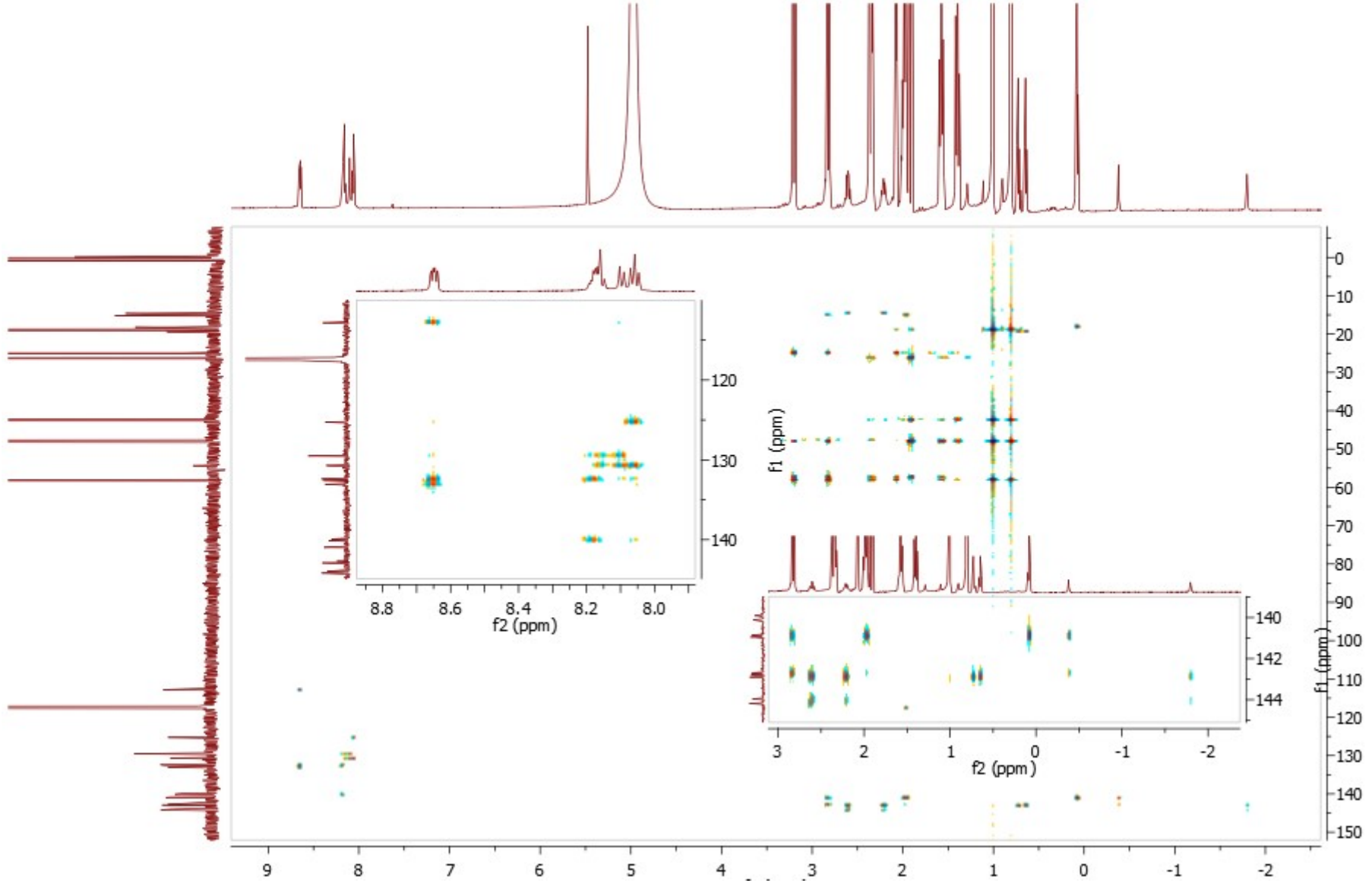

Figure S36. ^{1}H–^{13}C HMBC spectrum of α_4-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

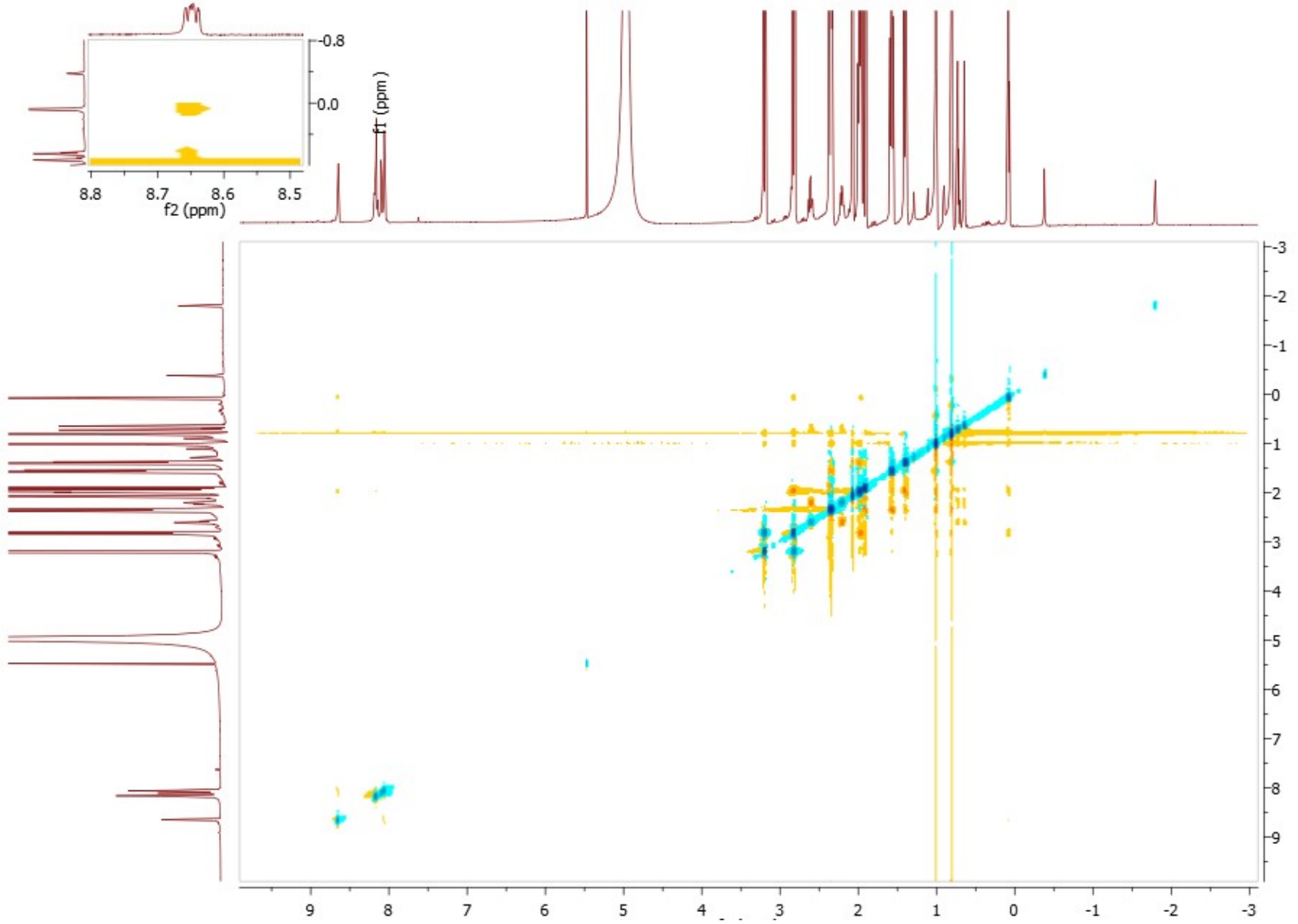
-0.8
0.0
f1 (ppm)
8.8
8.7
8.6
8.5
f2 (ppm)
-3
-2
-1
0
1
2
3
4
5
6
7
8
9
9
8
7
6
5
4
3
2
1
0
-1
-2
-3

Figure S37. ^{1}H–^{1}H ROESY spectrum of α_4-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

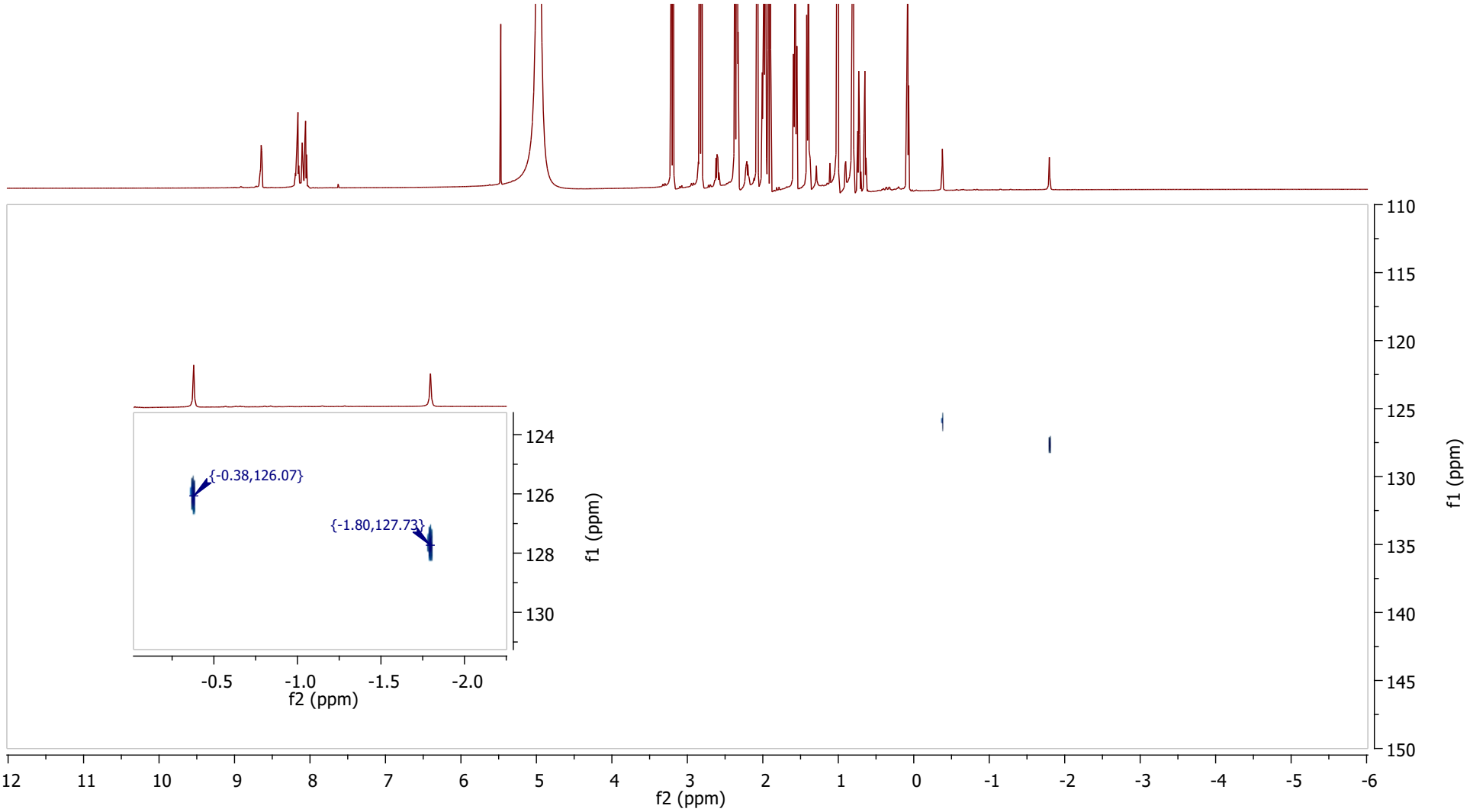

Figure S38. ^{1}H–^{15}N HSQC spectrum of α_4-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

α_4-P·10CSA(SR)

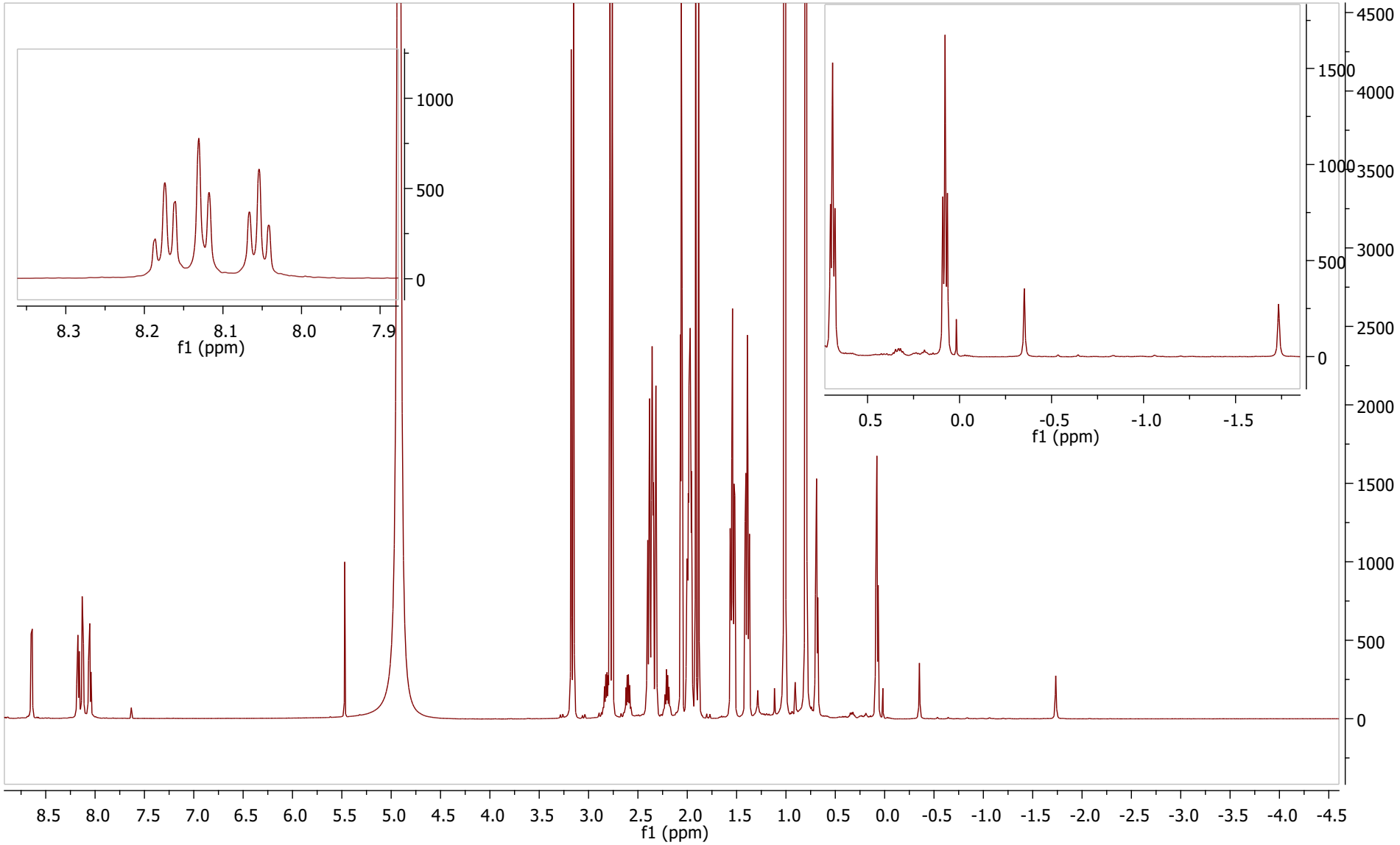

Figure S39. ^{1}H NMR spectrum of **α_4-P·10CSA(SR)** with expansion of areas of interest (600 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

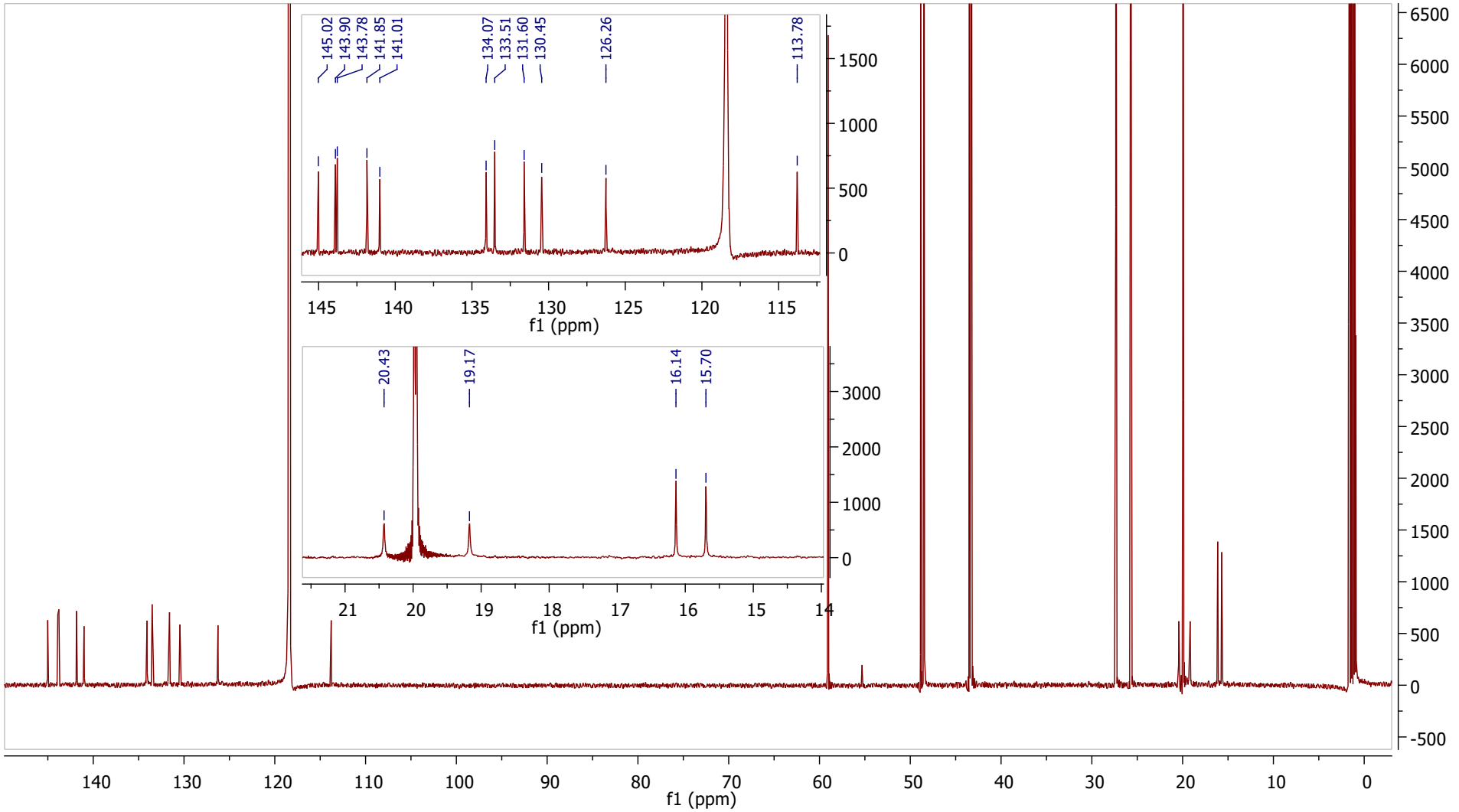

Figure S40. ^{13}C NMR spectrum of α_4-**P·10CSA(SR)** with expansion of areas of interest (151 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

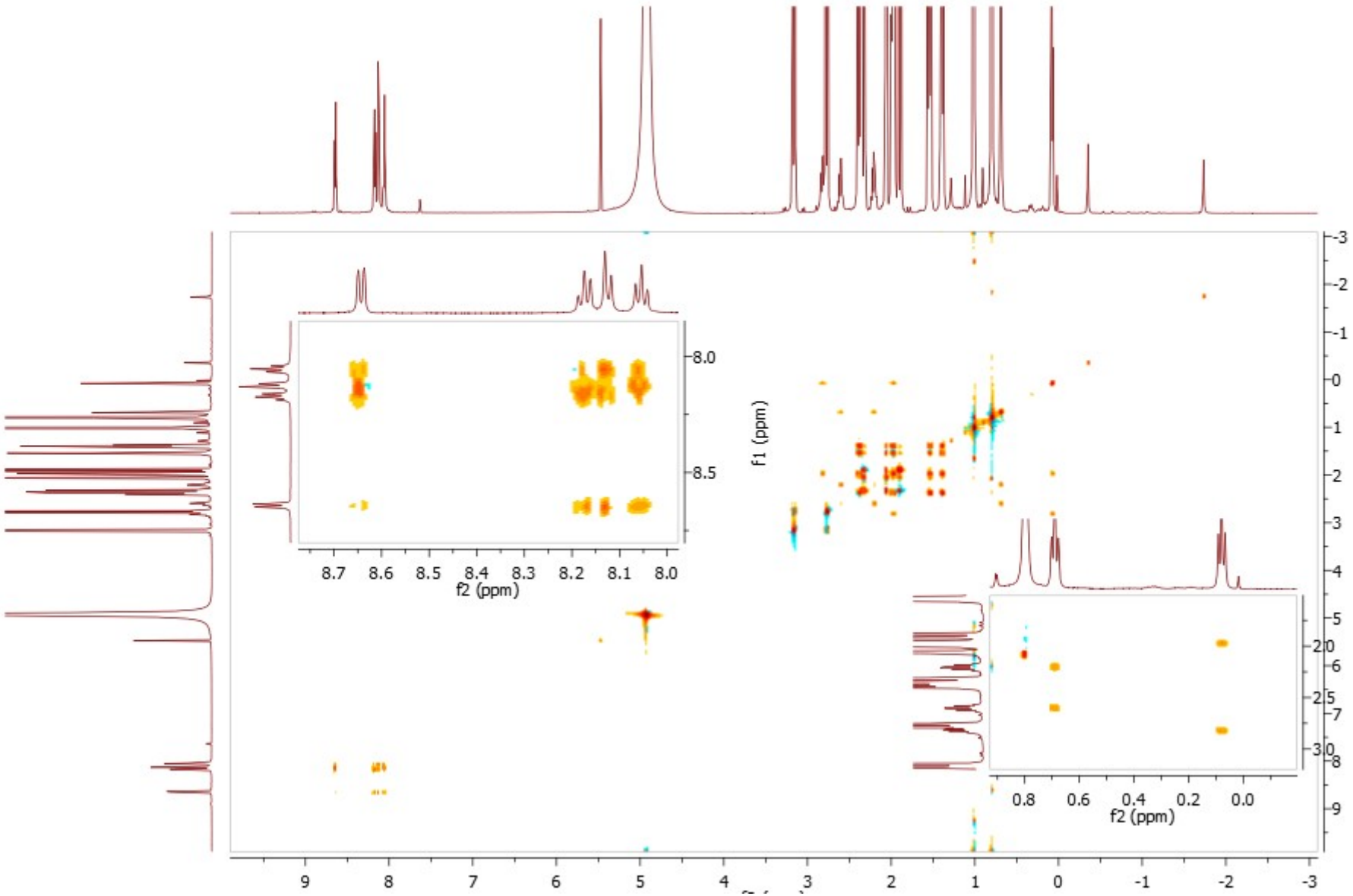

Figure S41. $^{1}H–^{1}H$ TOCSY spectrum of α_4-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

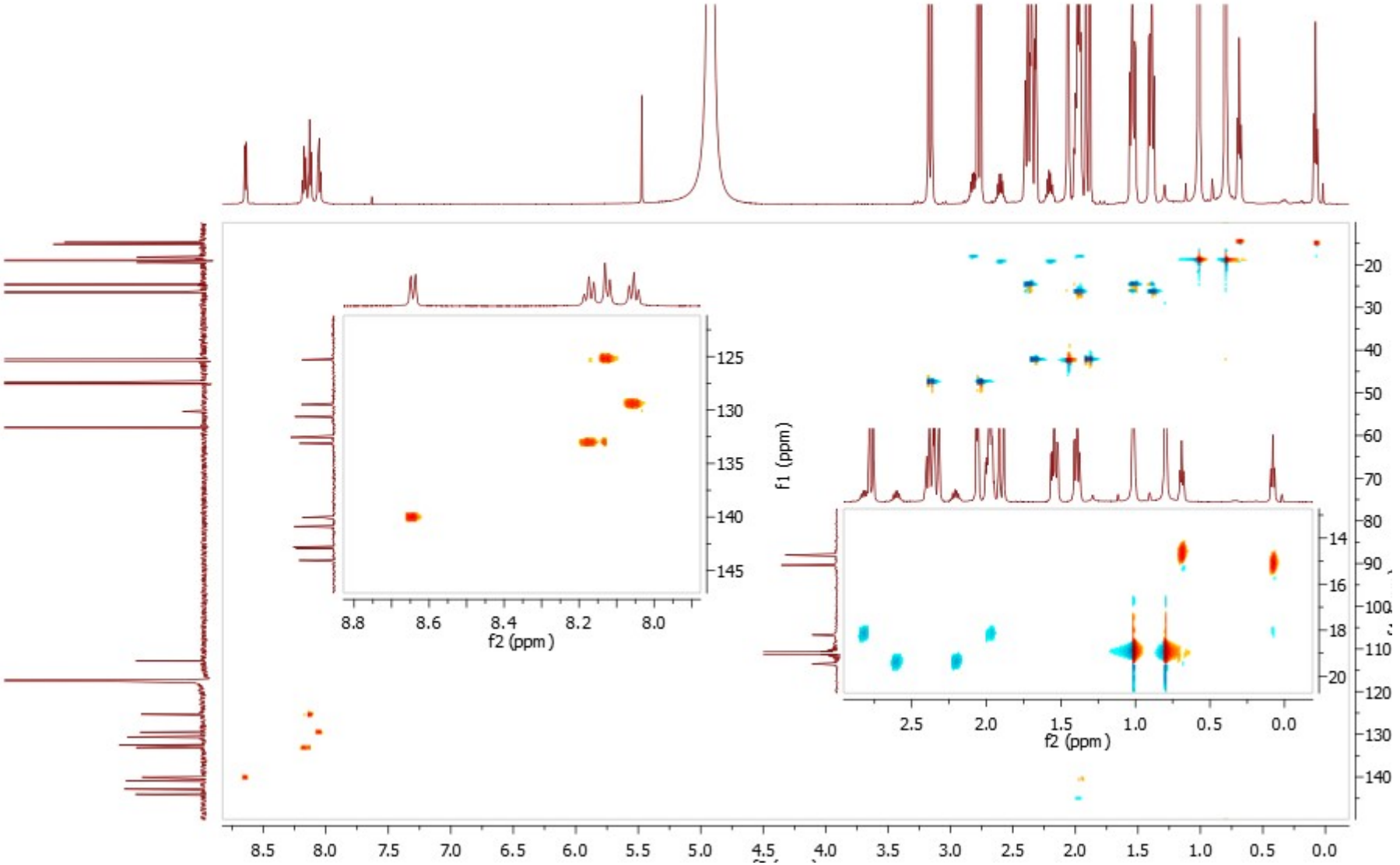

Figure S42. $^{1}H–^{13}C$ HSQC spectrum of α_4-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

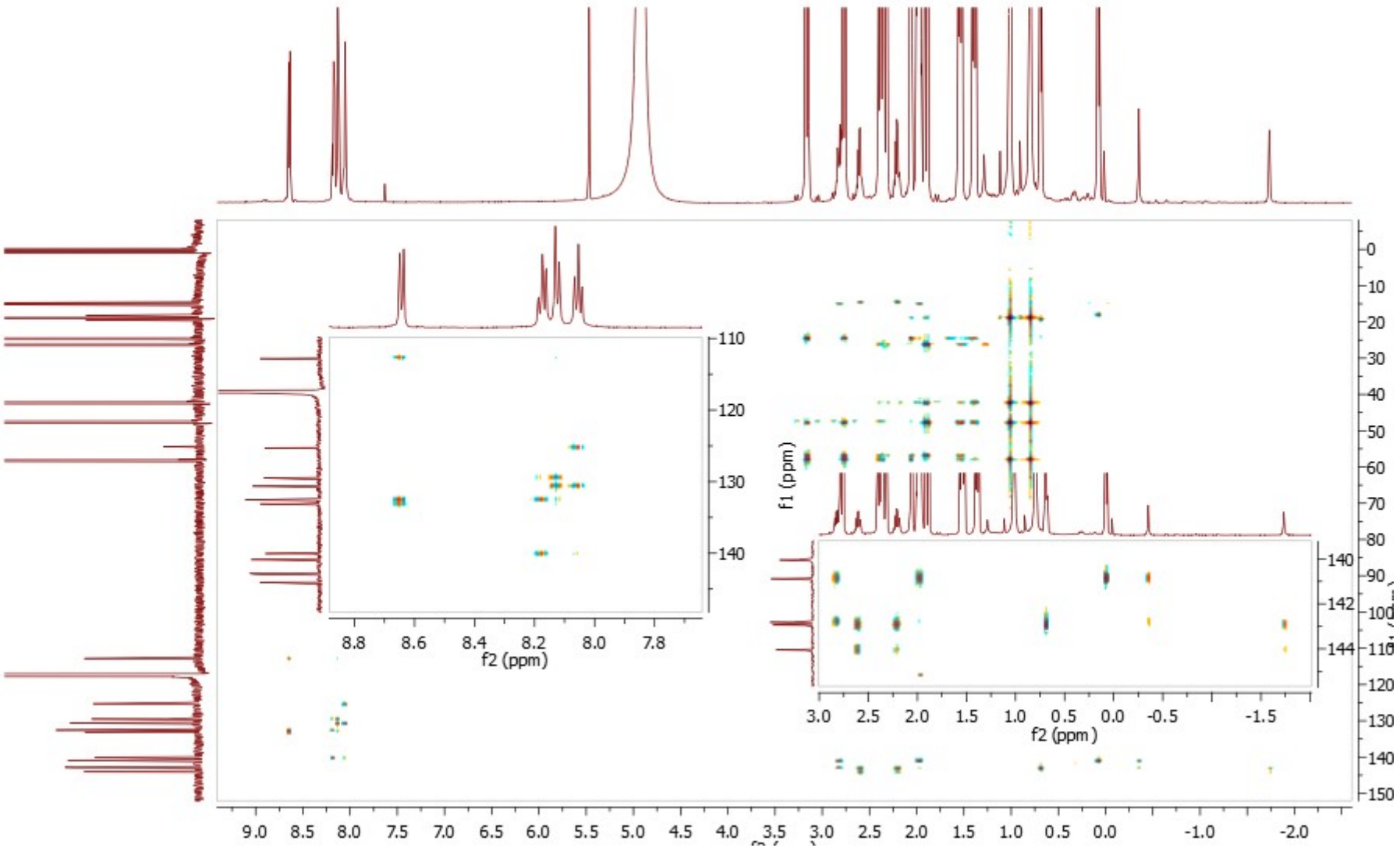

Figure S43. $^{1}H–^{13}C$ HMBC spectrum of α_4-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

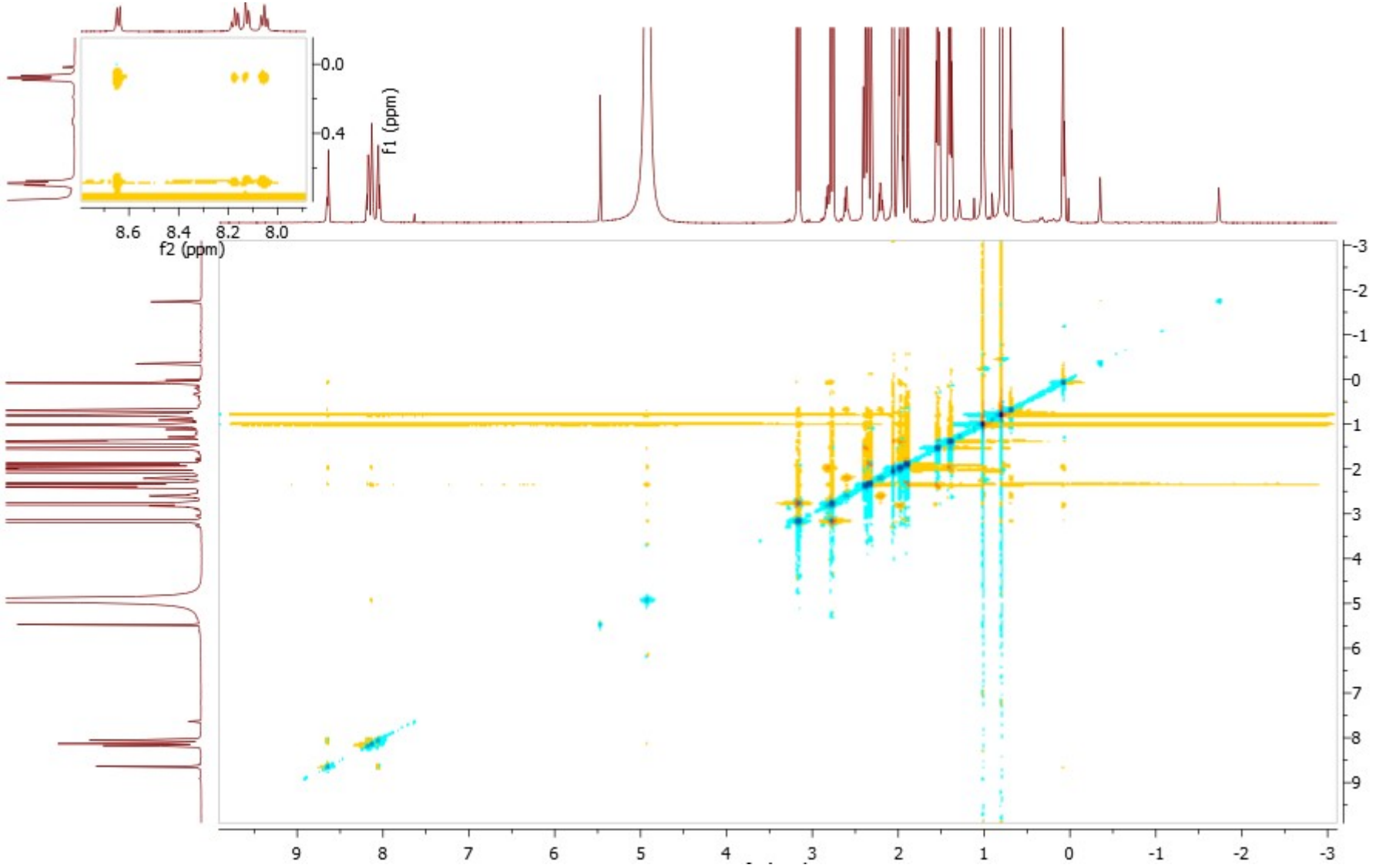

Figure S44. ^{1}H–^{1}H ROESY spectrum of α_4-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

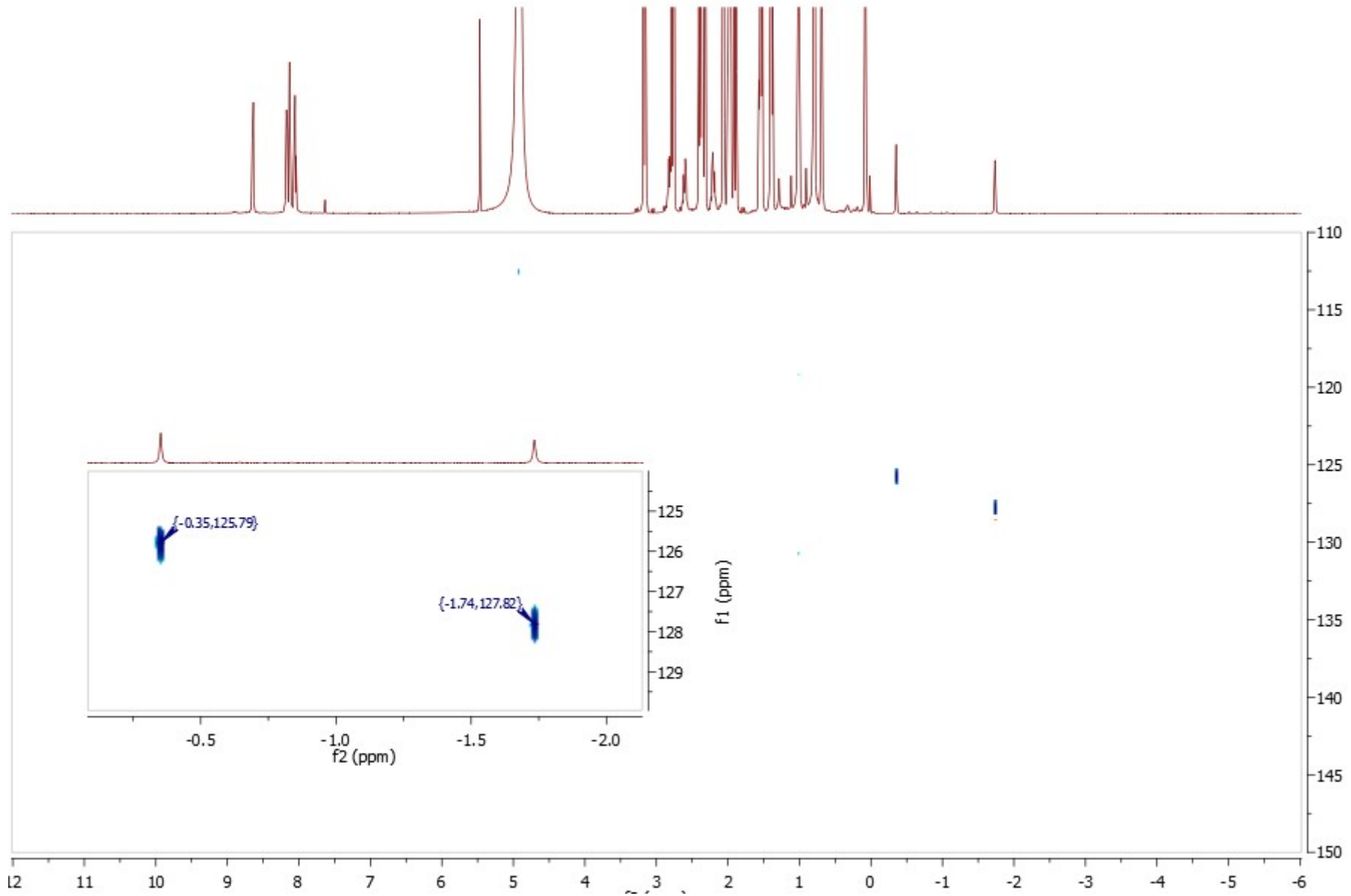

Figure S45. $^{1}H–^{15}N$ HSQC spectrum of α_4-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C)

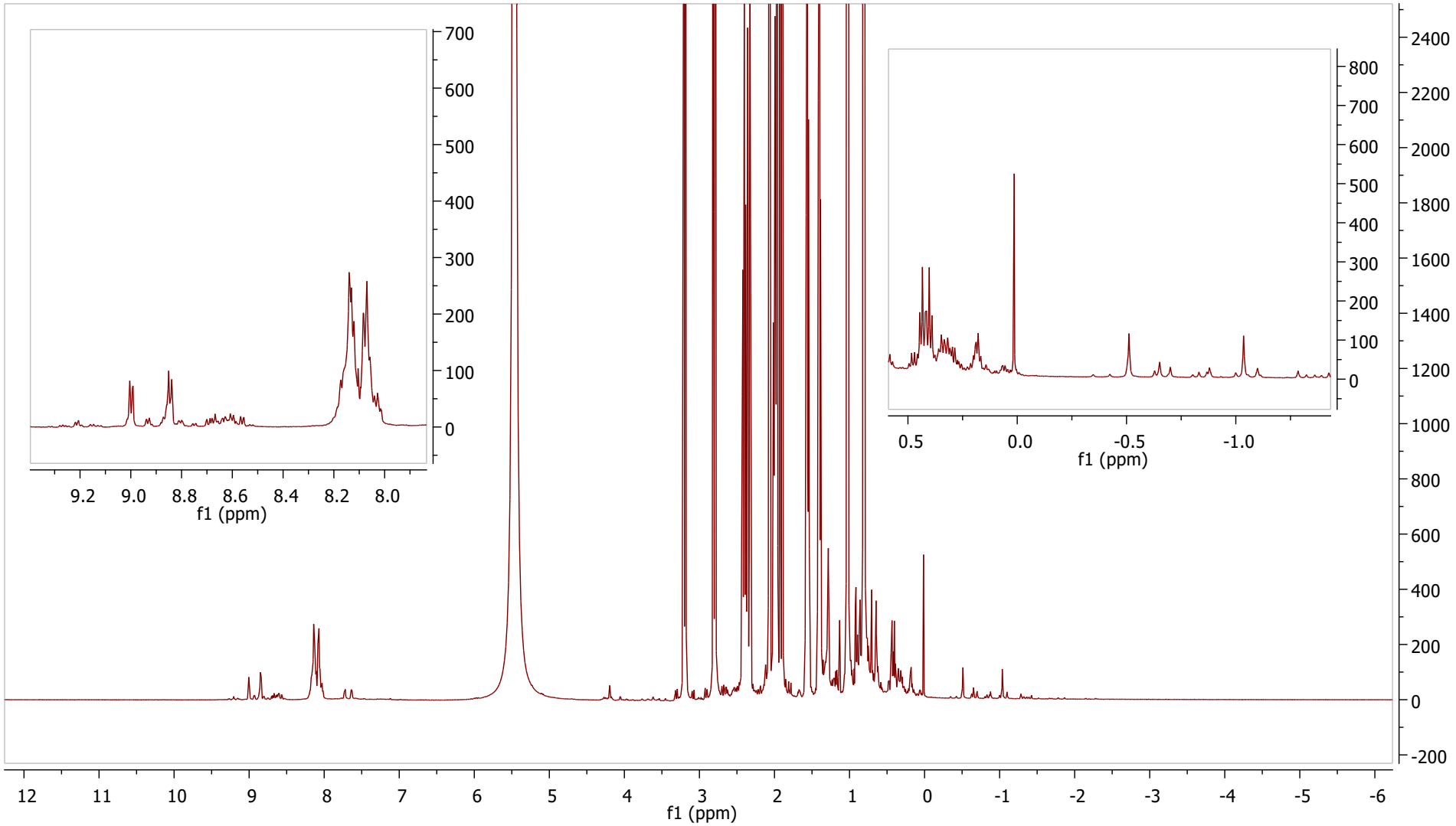

Figure S46. ^{1}H NMR spectrum of *αβαβ*-**P·10CSA(S)** with the expansion of areas of interest (600 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

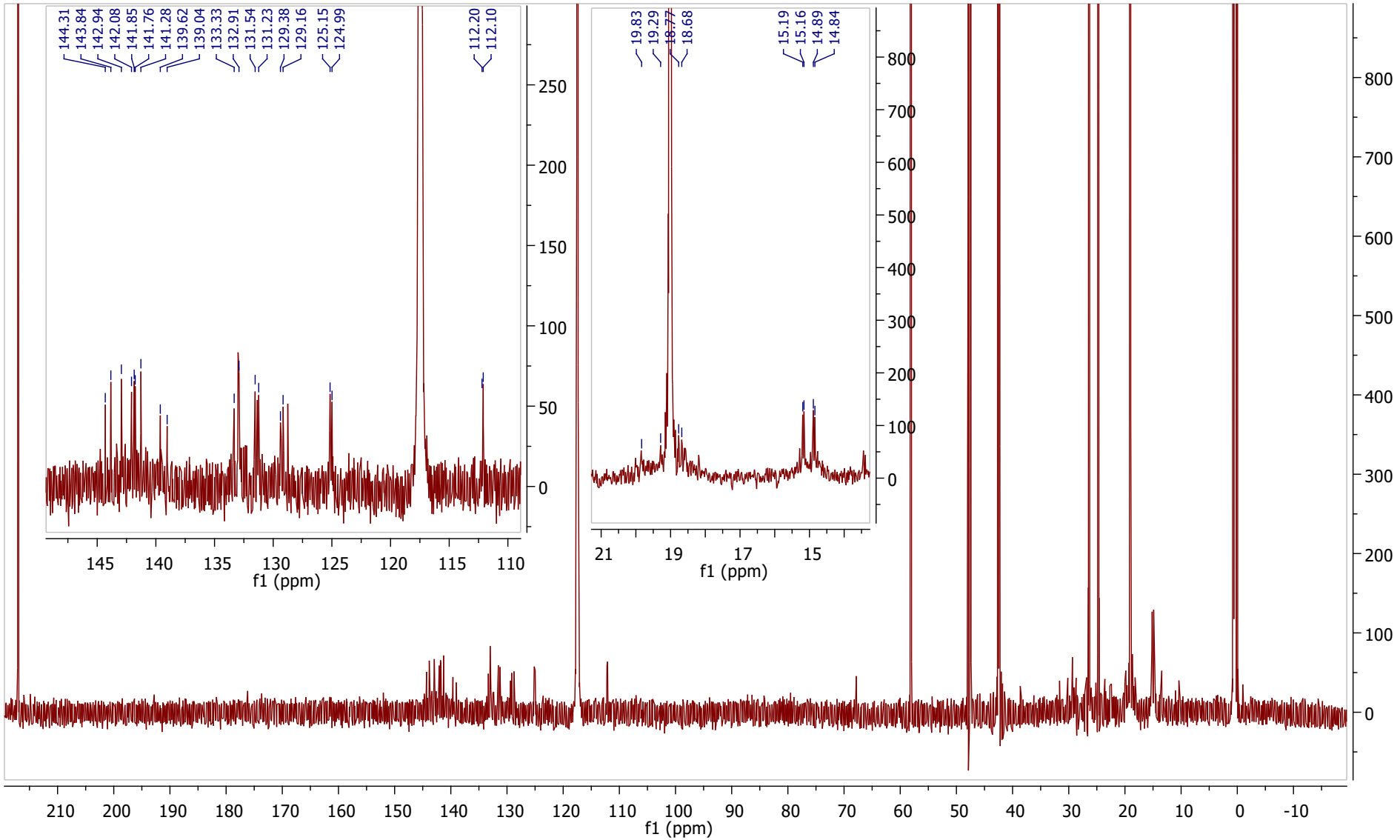

Figure S47. ^{13}C NMR spectrum of ***αβαβ*-P·10CSA(S)** with expansion of areas of interest (151 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

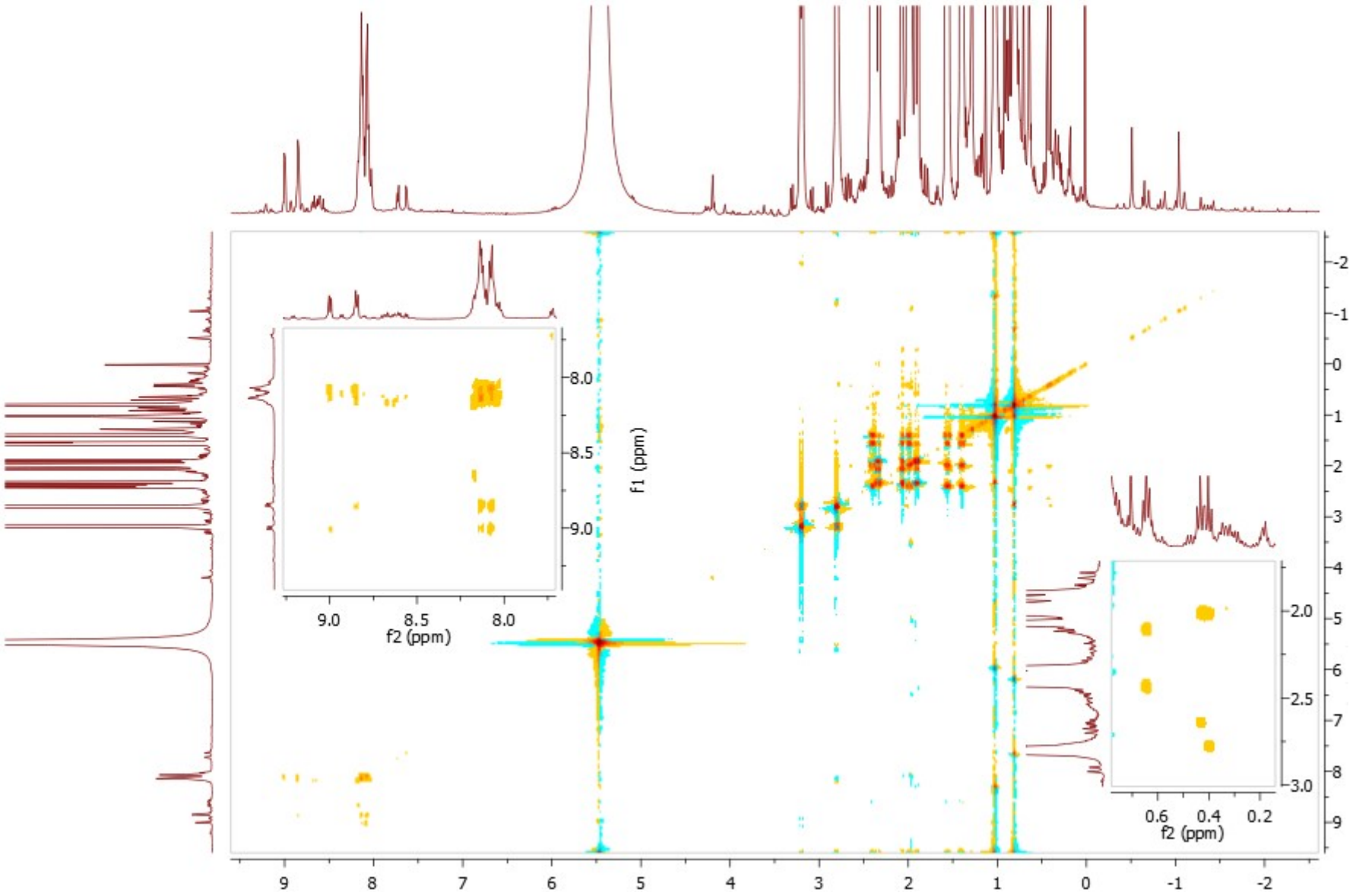

Figure S48. $^{1}H–^{1}H$ TOCSY spectrum of *αβαβ*-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

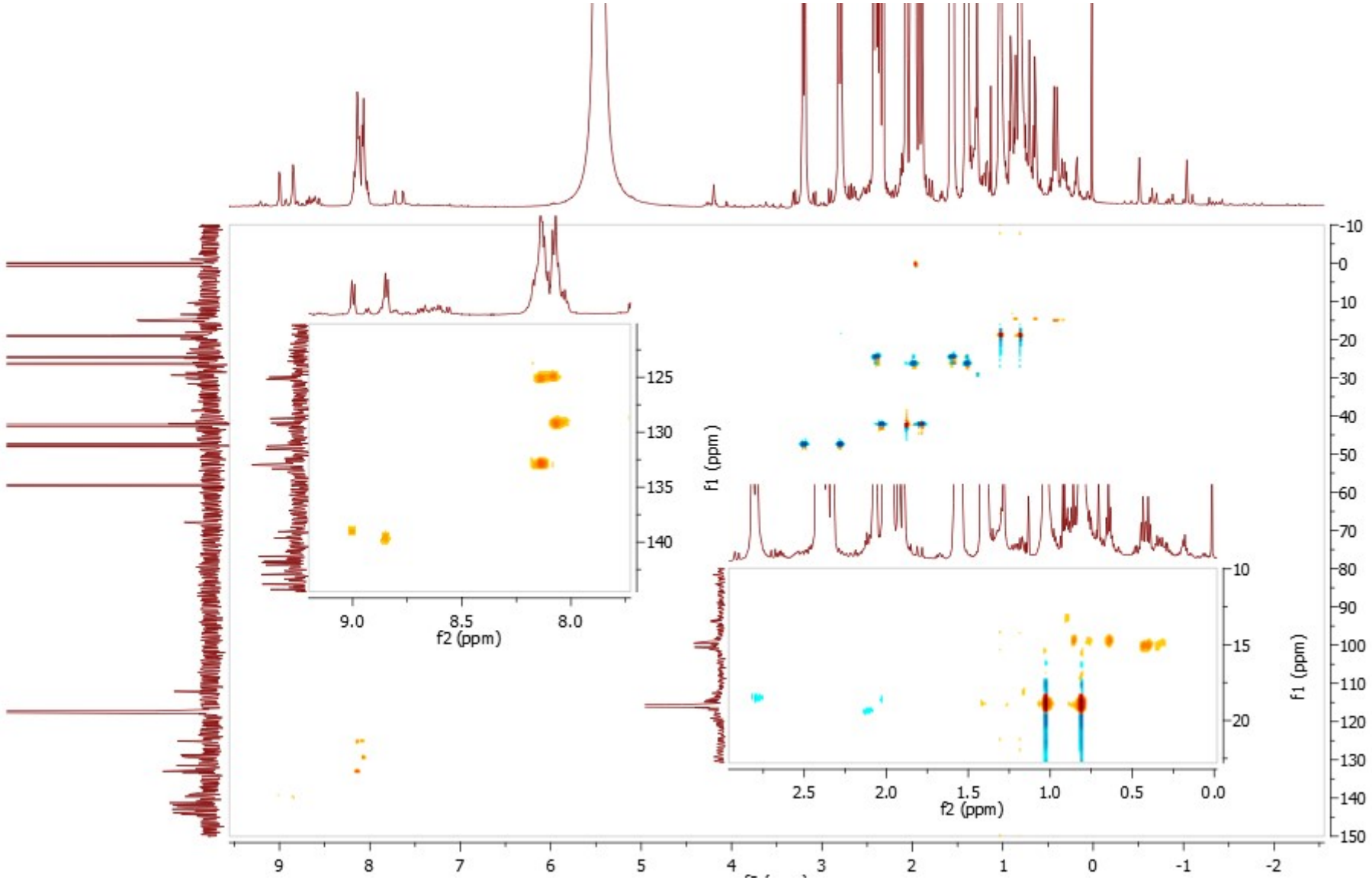

Figure S49. $^{1}H–^{13}C$ HSQC spectrum of *αβαβ*-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

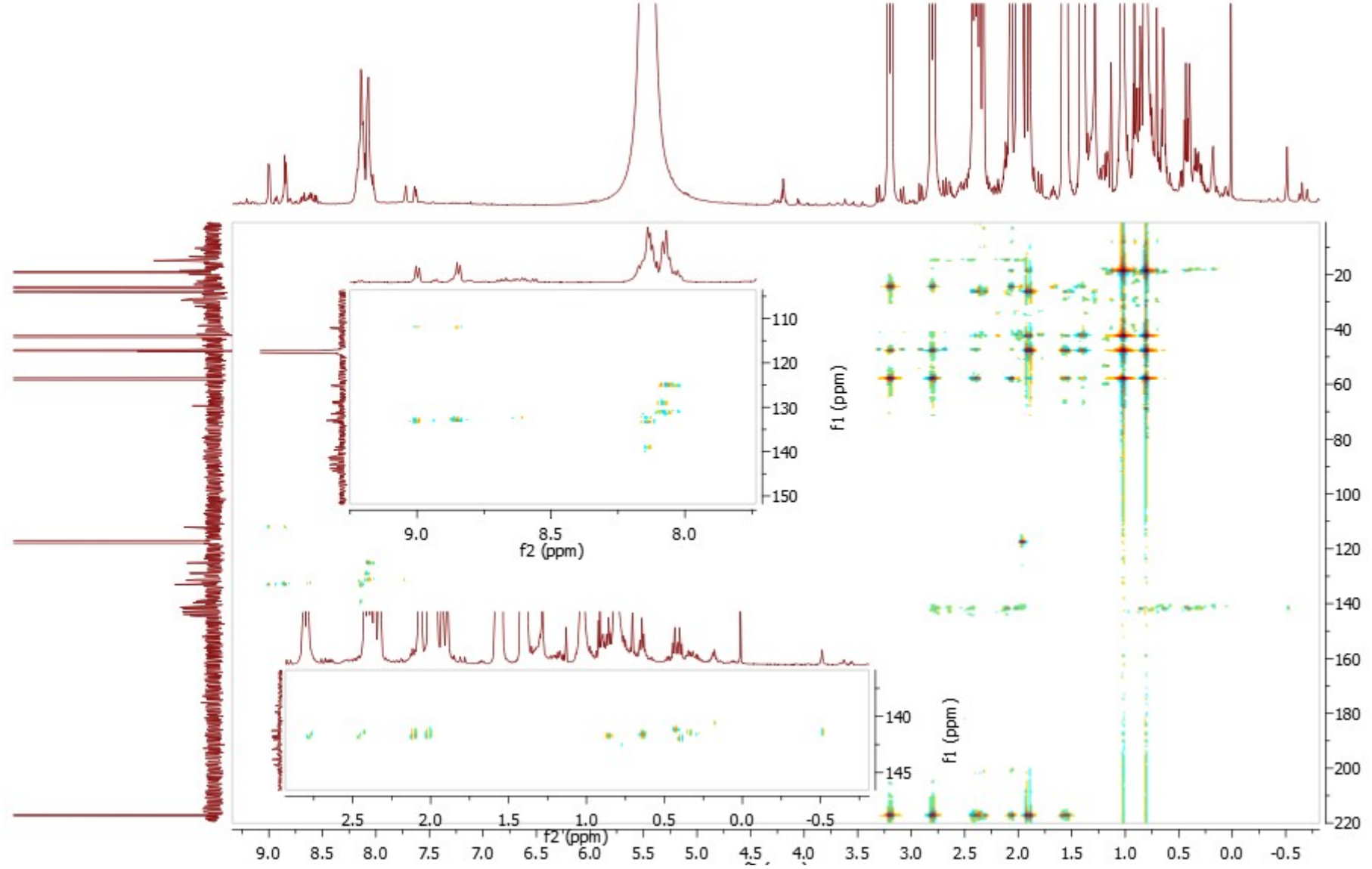

Figure S50. $^{1}H–^{13}C$ HMBC spectrum of *αβαβ*-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

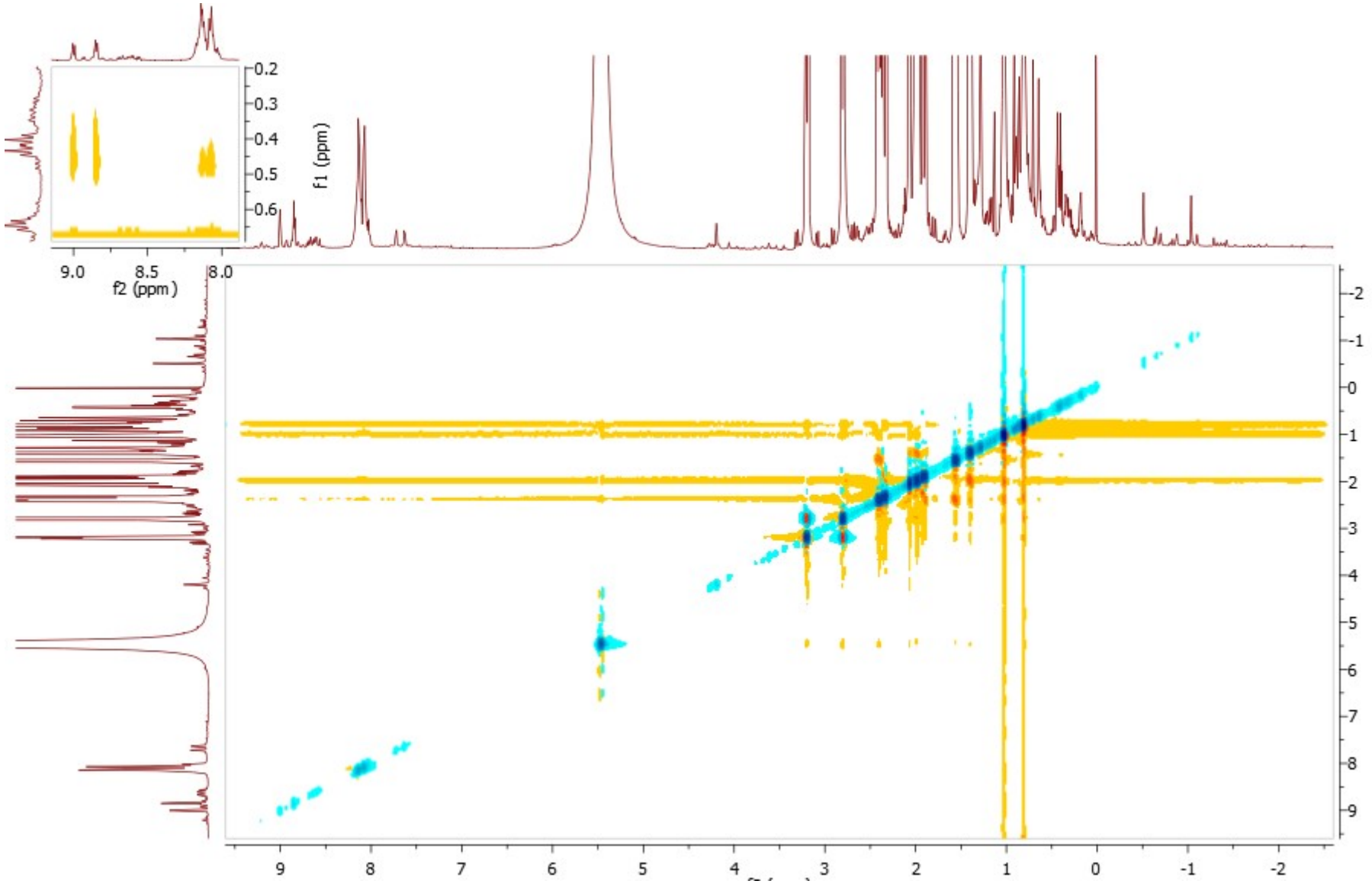

Figure S51. ^{1}H–^{1}H ROESY spectrum of *αβαβ*-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

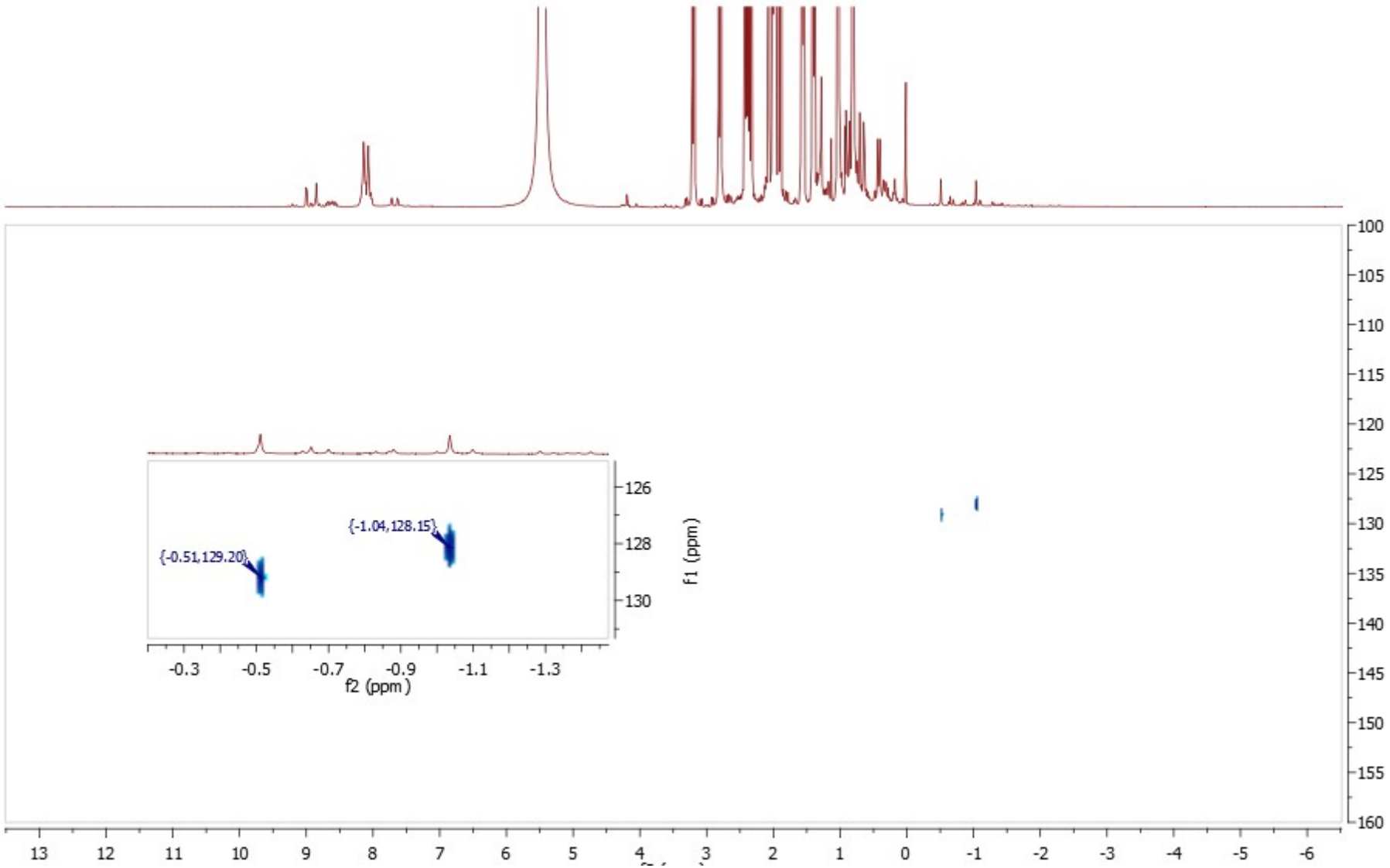

Figure S52. ^{1}H–^{15}N HSQC spectrum of *αβαβ*-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

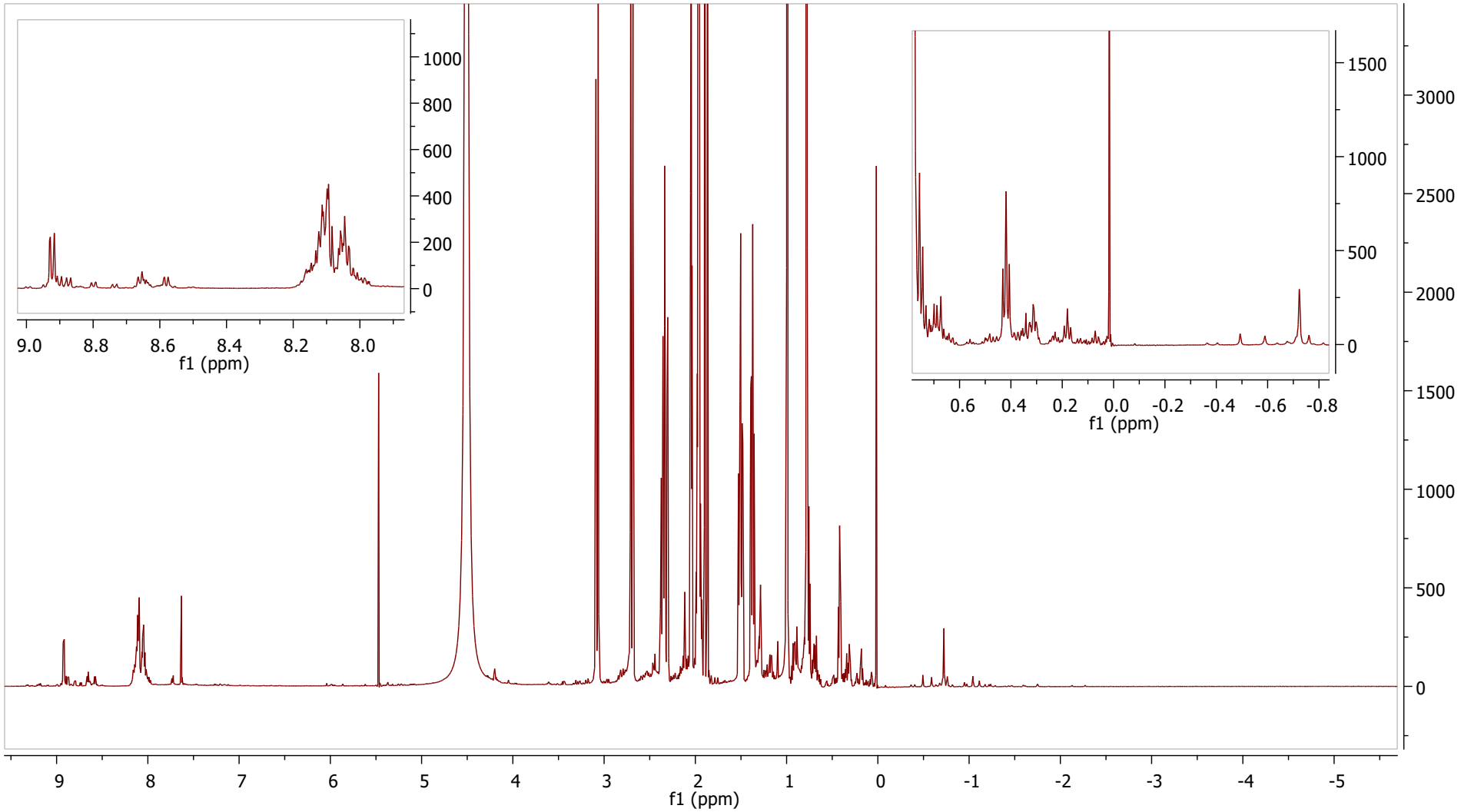

Figure S53. ^{1}H NMR spectrum of *αβαβ*-**P·10CSA(SR)** with expansion of areas of interest (600 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

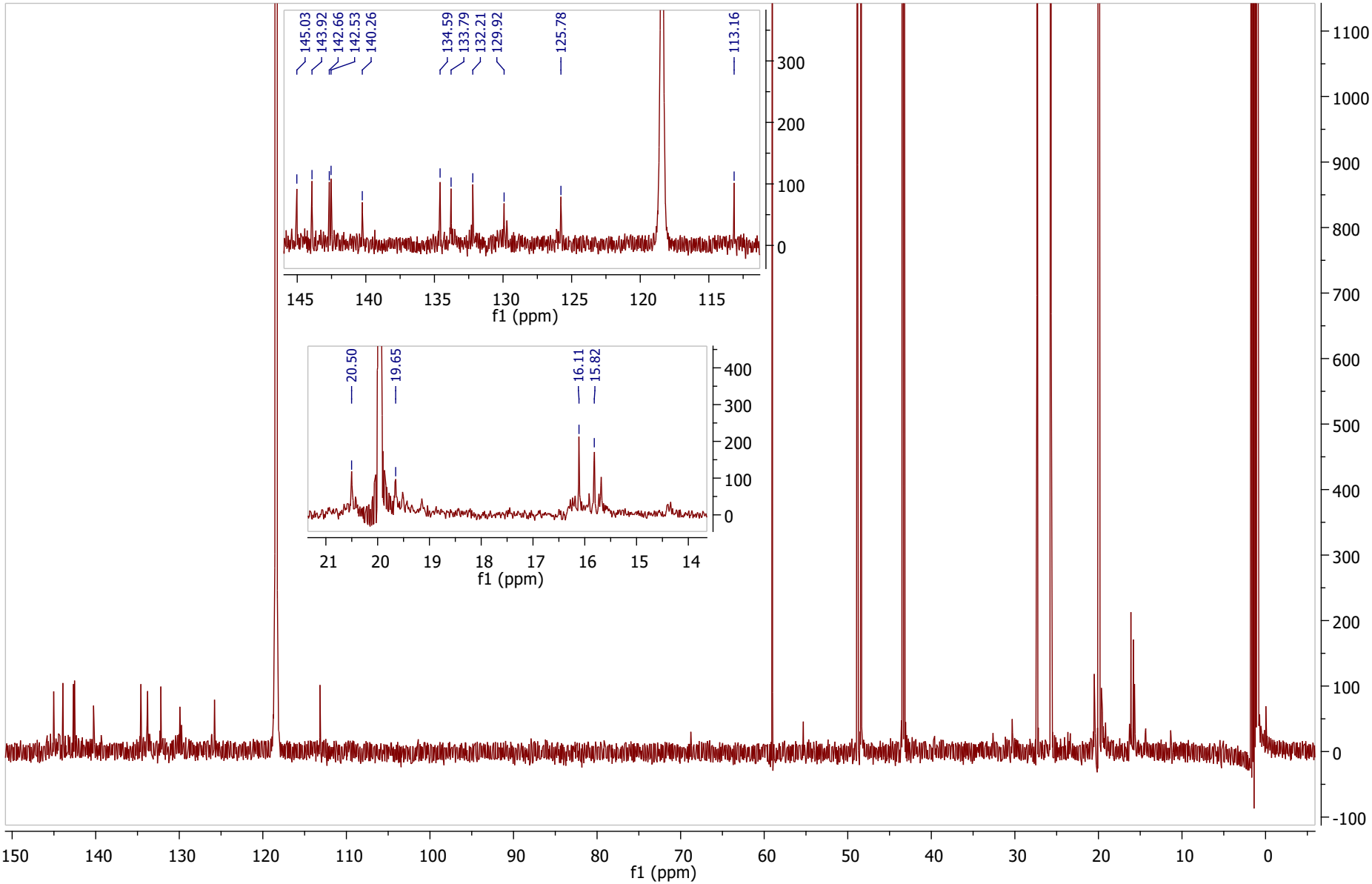

Figure S54. ^{13}C NMR spectrum of *αβαβ*-**P·10CSA(SR)** with expansion of areas of interest (151 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

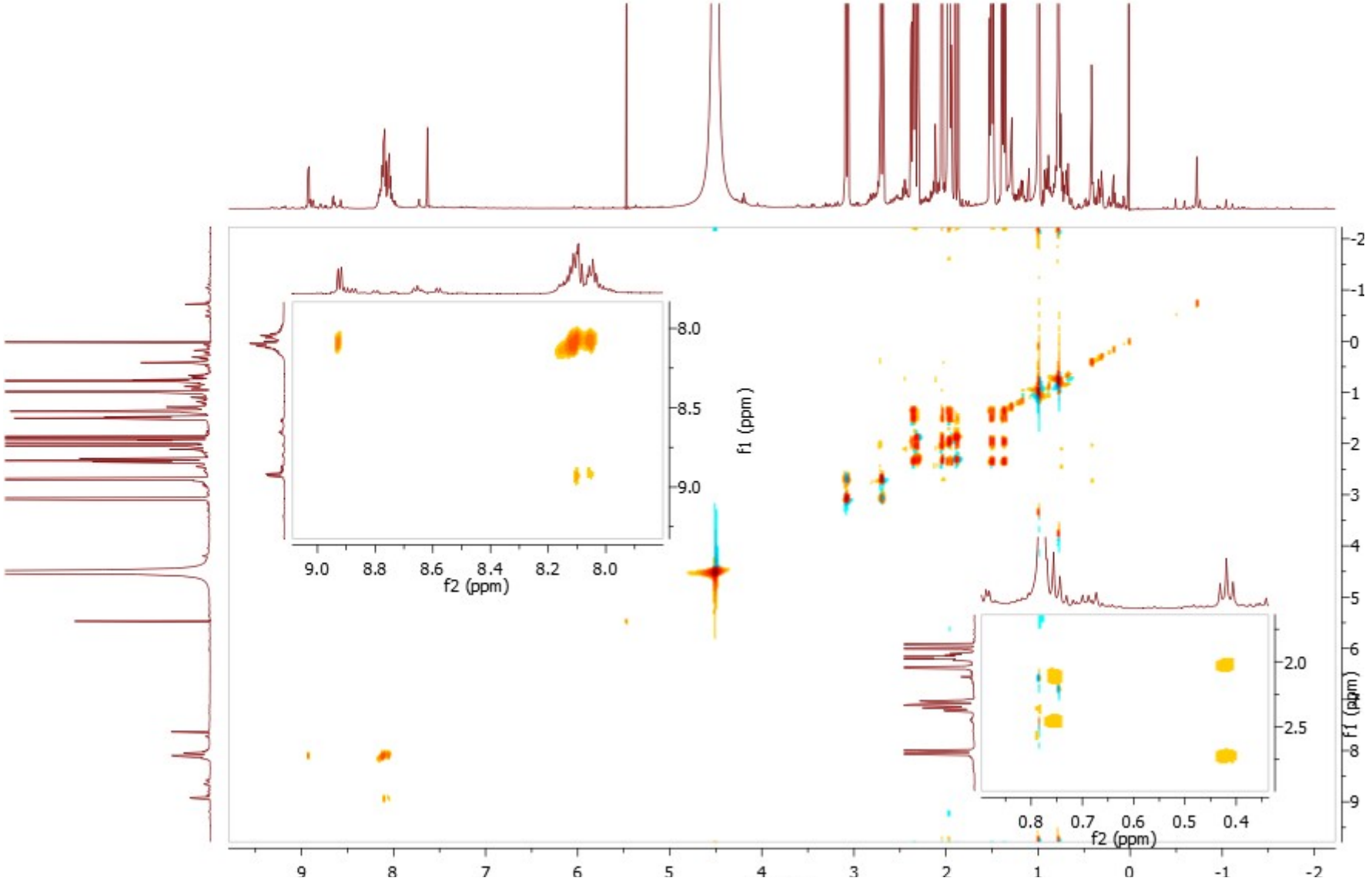

Figure S55. $^{1}H–^{1}H$ TOCSY spectrum of ***αβαβ*-P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

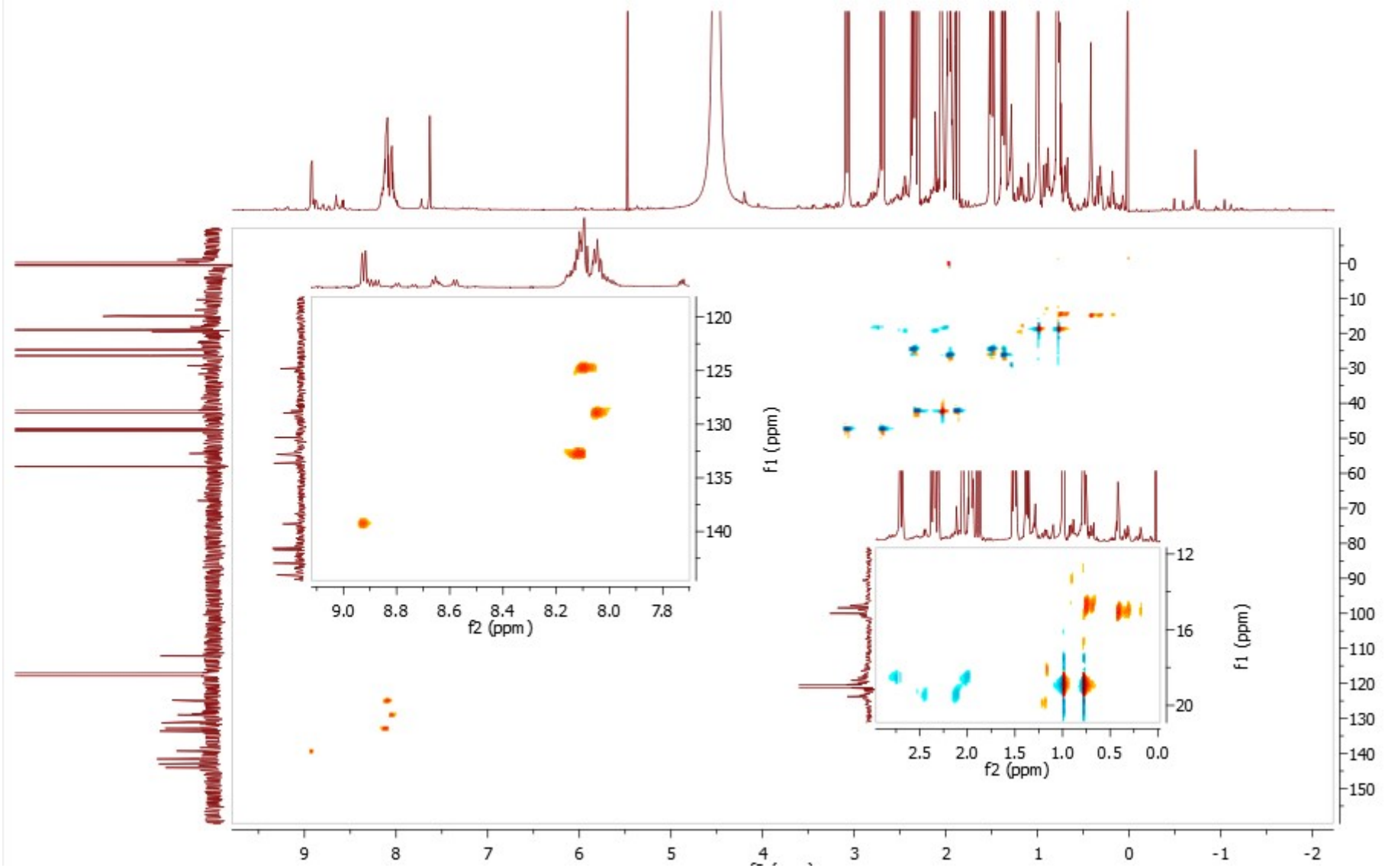

Figure S56. $^{1}H–^{13}C$ HSQC spectrum of *αβαβ*-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

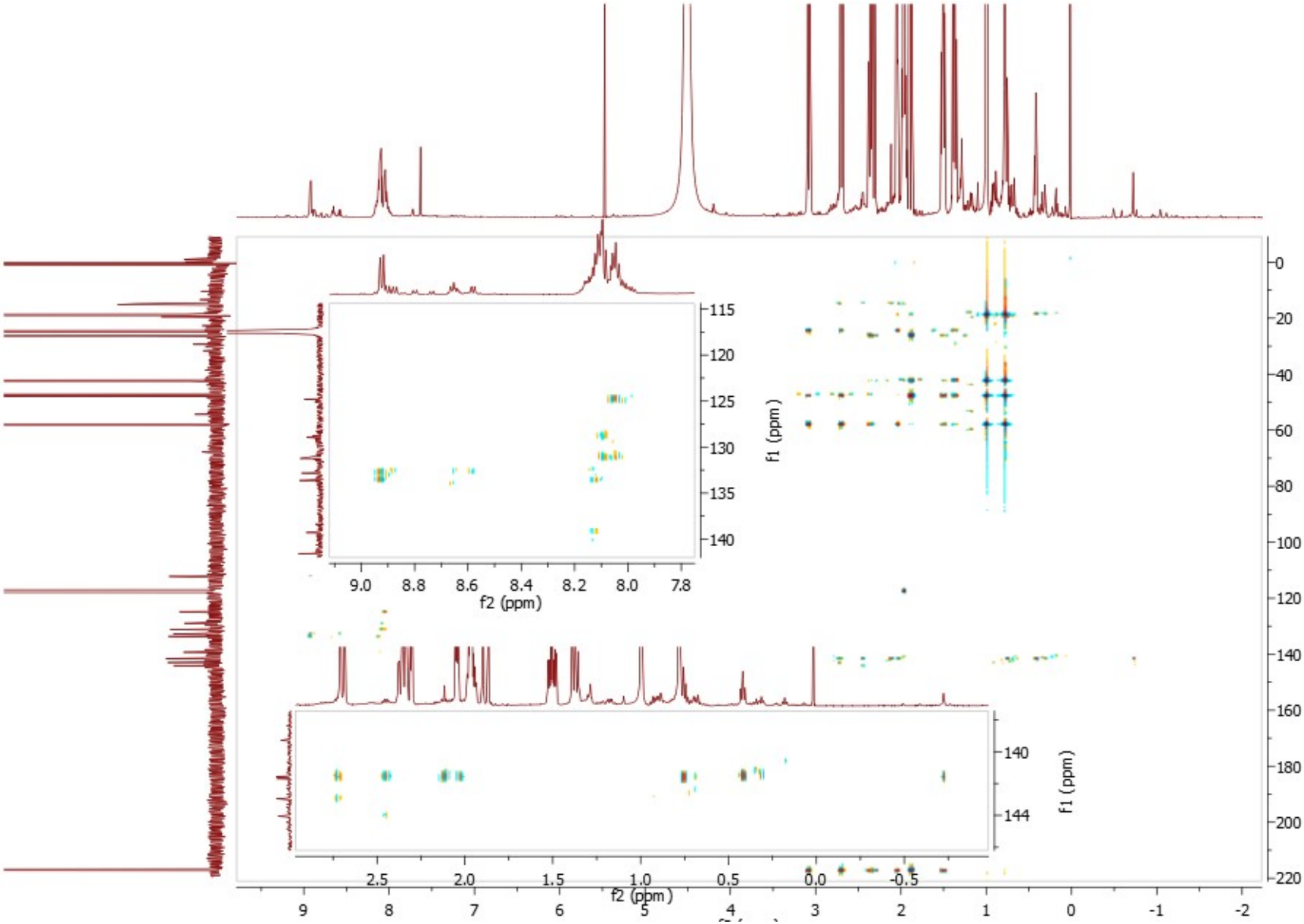

Figure S57. $^{1}H–^{13}C$ HMBC spectrum of *αβαβ*-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

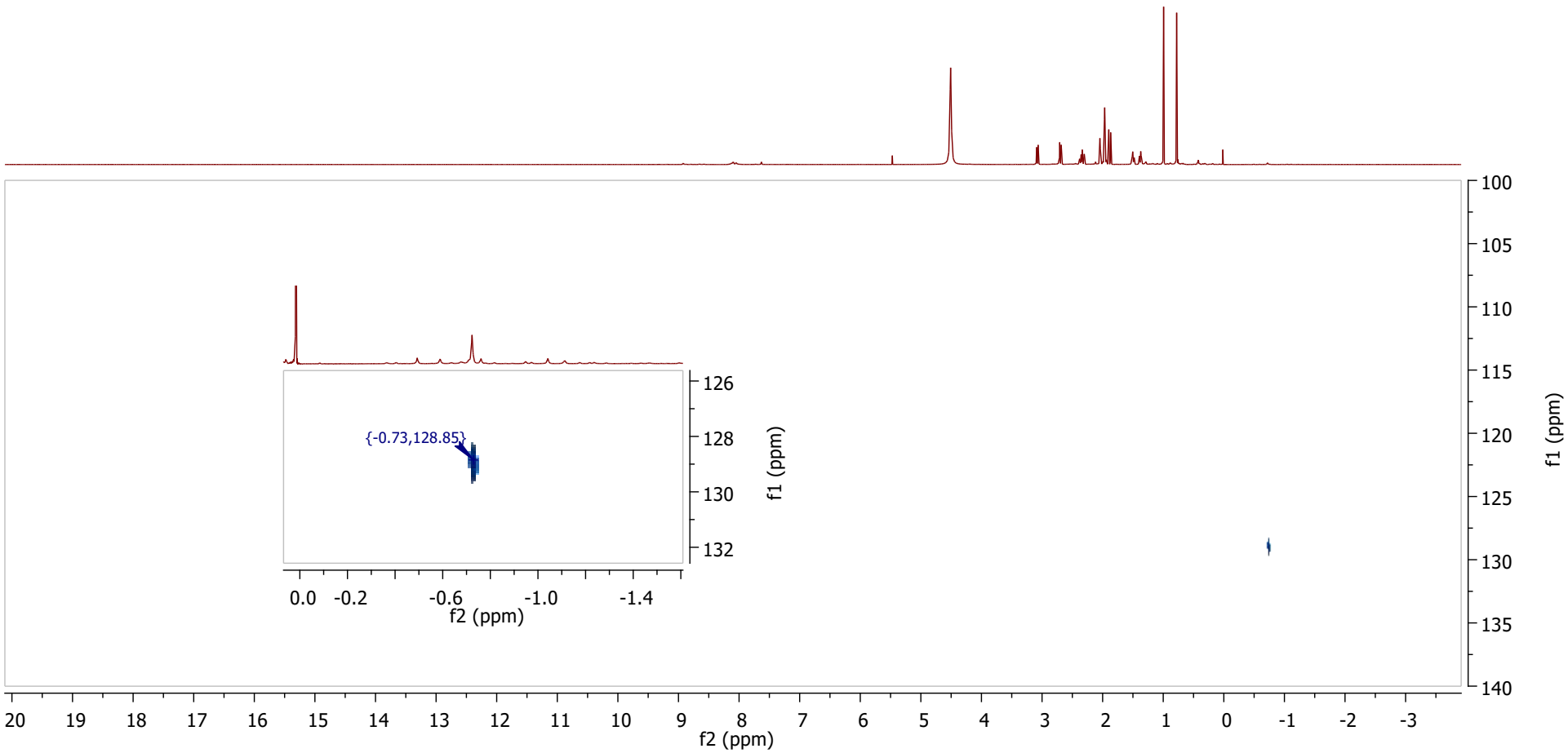

Figure S58. $^{1}H–^{15}N$ HSQC spectrum of *αβαβ*-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C)

$\alpha_3\beta$-P·10CSA(S)

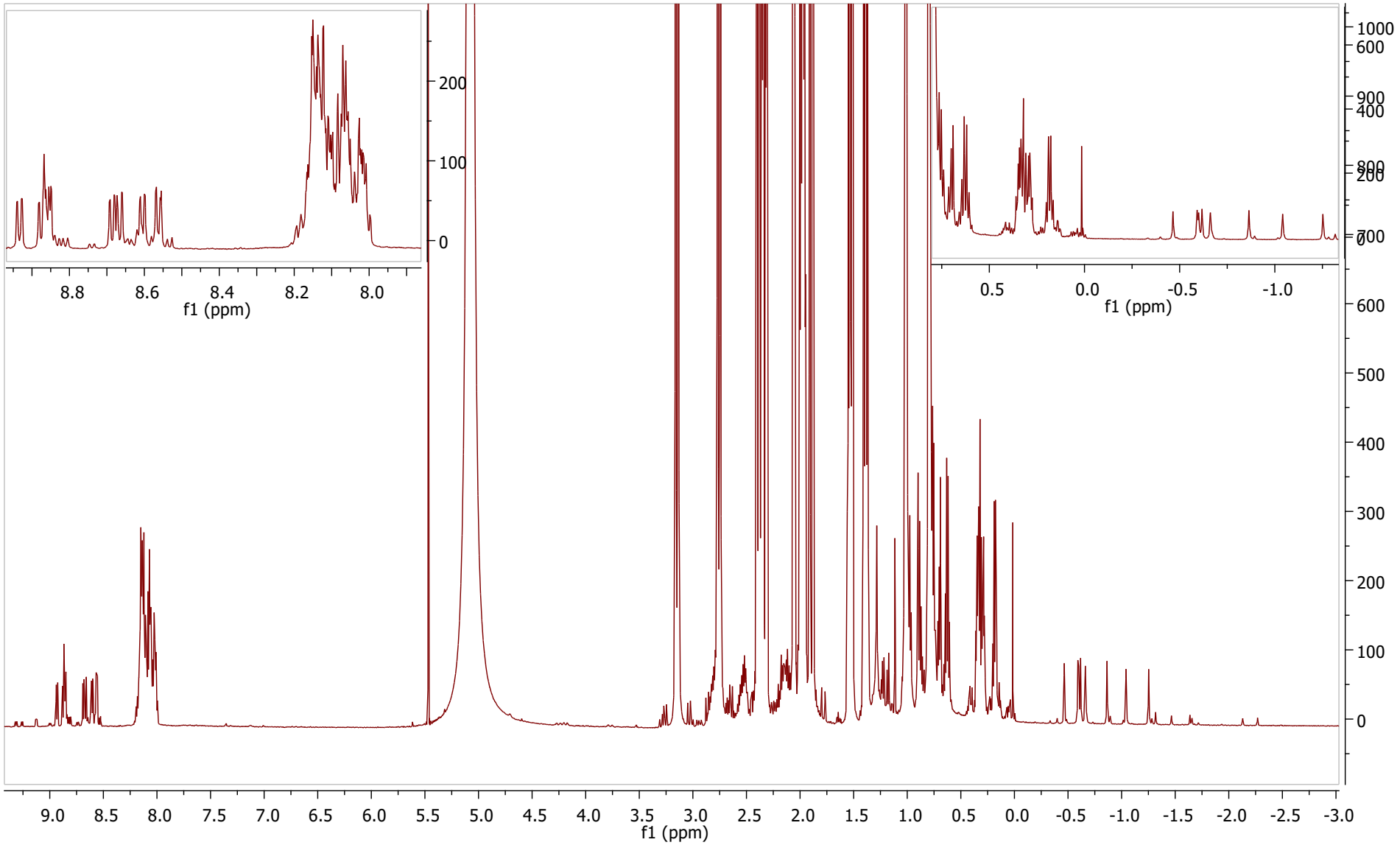

Figure S59. ^{1}H NMR spectrum of $\alpha_3\beta$-**P·10CSA(S)** with the expansion of areas of interest (600 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

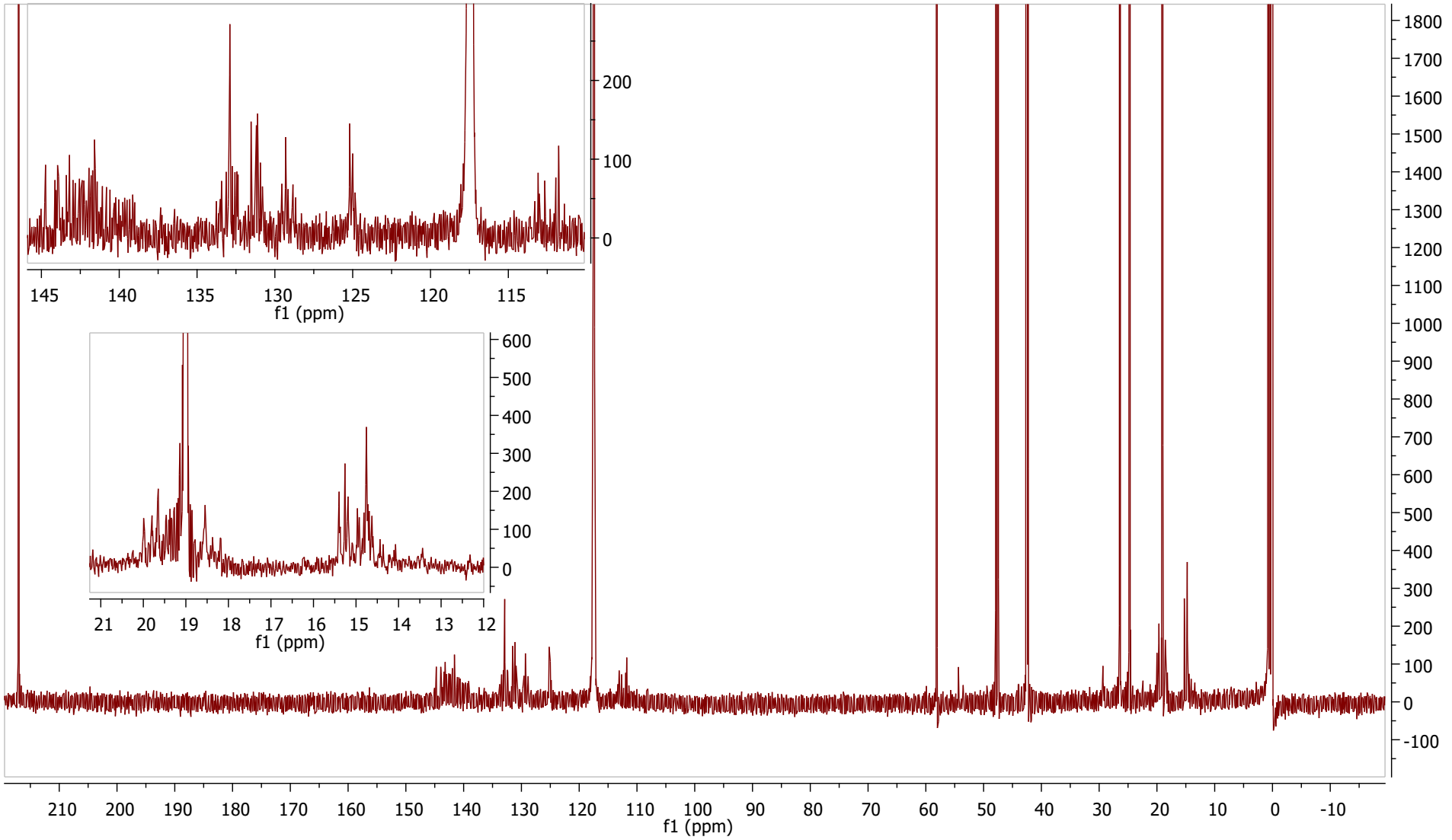

Figure S60. ^{13}C NMR spectrum of $\alpha_3\beta$-**P·10CSA(S)** with expansion of areas of interest (151 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

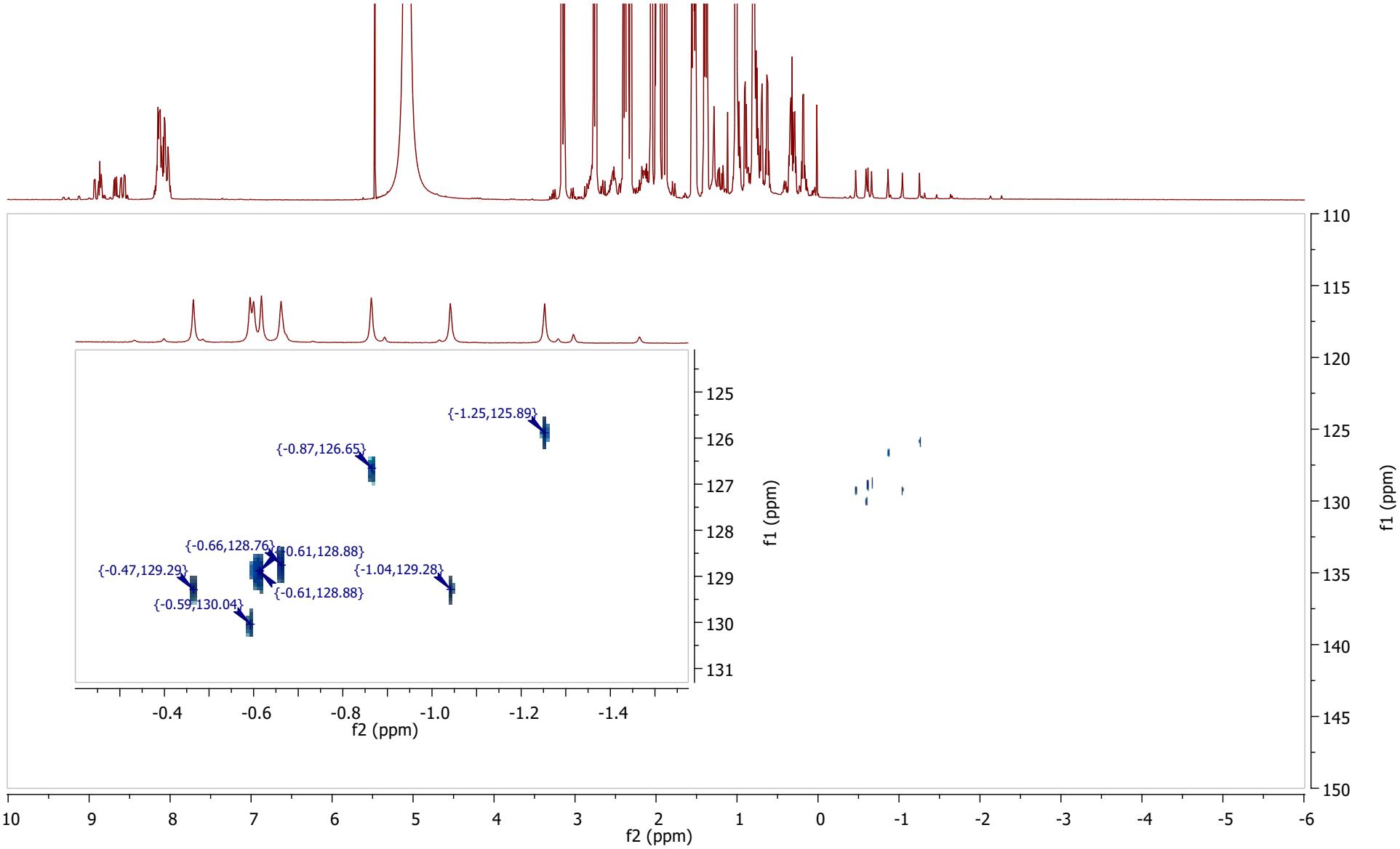

Figure S61. $^{1}H–^{15}N$ HSQC spectrum of $\alpha_3\beta$-**P·10CSA(S)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S), 25 °C).

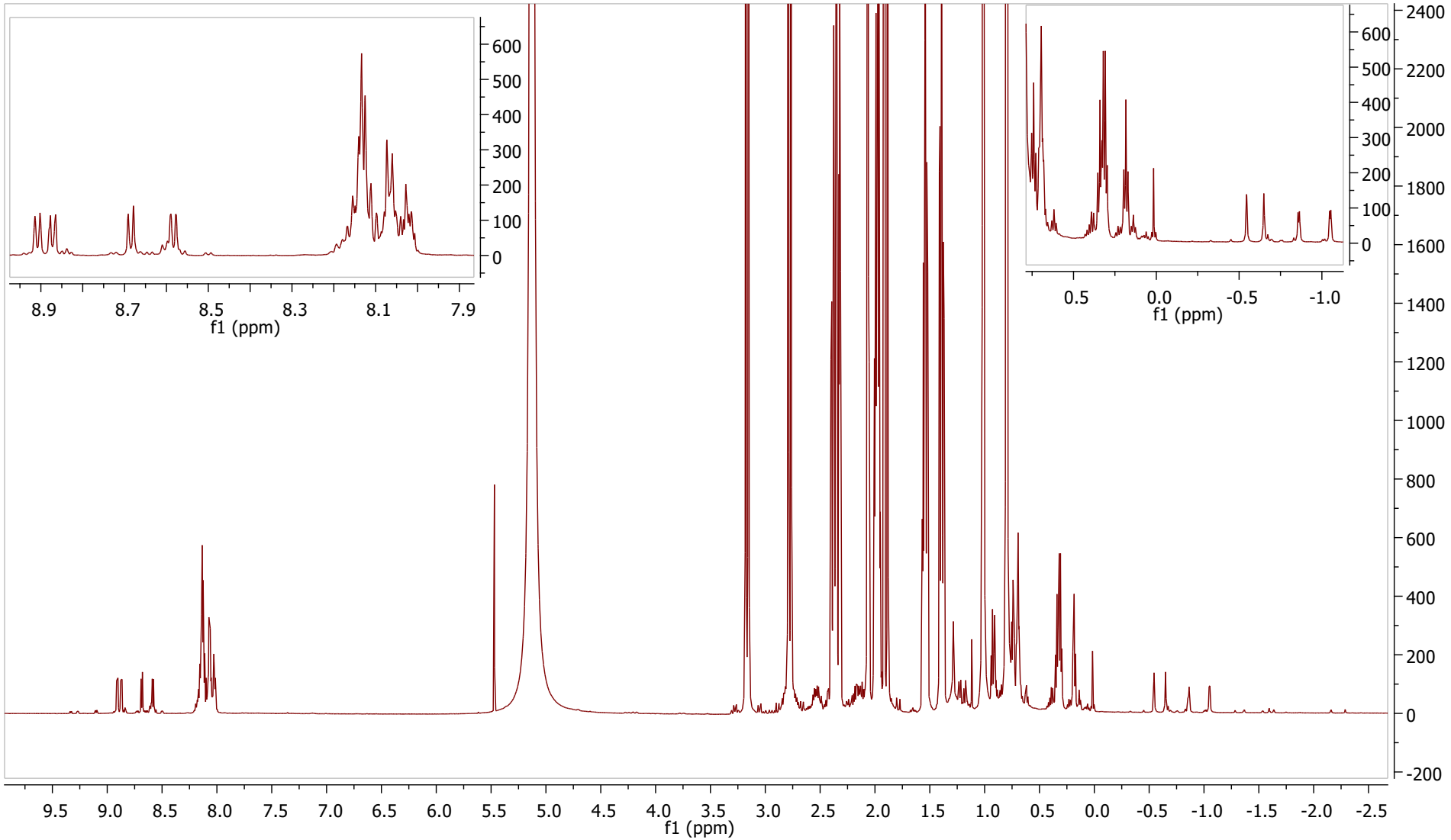

Figure S62. ^{1}H NMR spectrum of $\alpha_3\beta$-**P·10CSA(SR)** with expansion of areas of interest (600 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

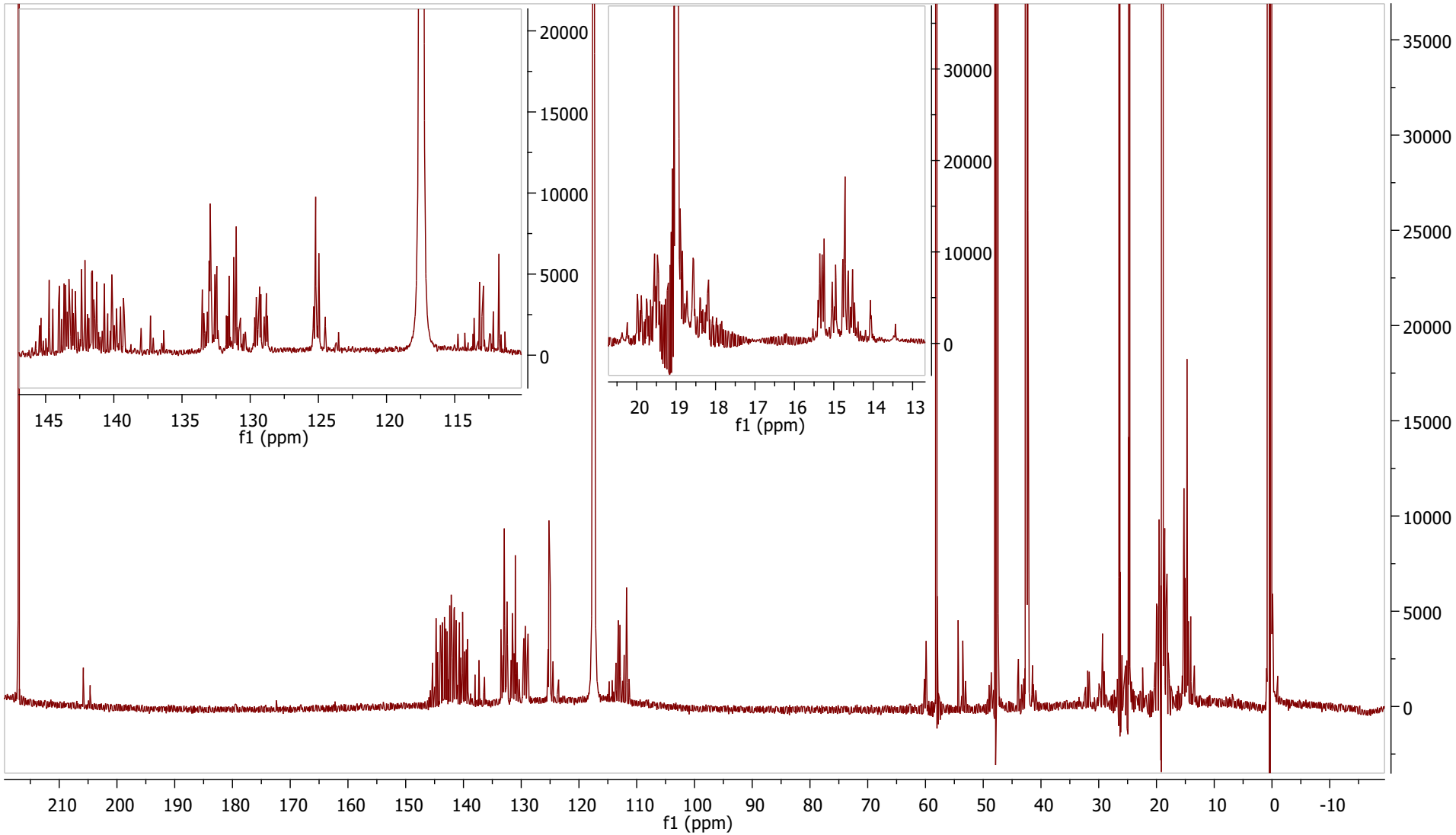

Figure S63. ^{13}C NMR spectrum of $\alpha_3\beta$-**P·10CSA(SR)** with expansion of areas of interest (151 MHz, acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C).

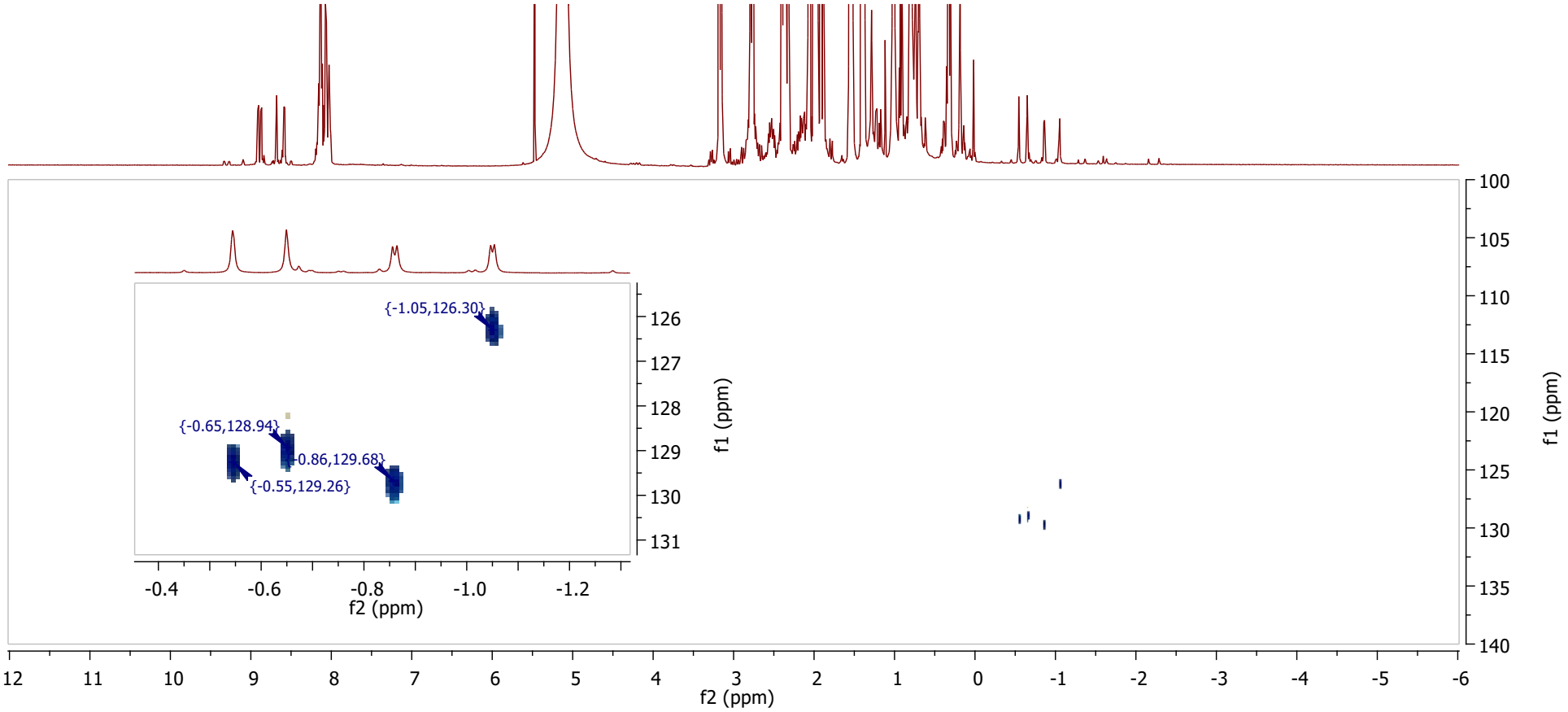

Figure S64. $^{1}H–^{15}N$ HSQC spectrum of $\alpha_3\beta$-**P·10CSA(SR)** with expansion of areas of interest (acetonitrile-d_3, 20 eq. of 10CSA(S) and 10CSA(R), 25 °C)

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2. C. J. Kingsbury and M. O. Senge, *Coord. Chem. Rev.*, 2021, **431**, 213760–213783.
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