



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 12, 2026 – 09:01 PM UTC

PDB ID : 1L0L / pdb\_00001l0l  
Title : structure of bovine mitochondrial cytochrome bc1 complex with a bound fungicide famoxadone  
Authors : Gao, X.; Wen, X.; Yu, C.A.; Esser, L.; Tsao, S.; Quinn, B.; Zhang, L.; Yu, L.; Xia, D.  
Deposited on : 2002-02-11  
Resolution : 2.35 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

|                                |   |                                                                  |
|--------------------------------|---|------------------------------------------------------------------|
| MolProbity                     | : | 4-5-2 with Phenix2.0                                             |
| Mogul                          | : | 2022.3.0, CSD as543be (2022)                                     |
| Xtriage (Phenix)               | : | NOT EXECUTED                                                     |
| EDS                            | : | NOT EXECUTED                                                     |
| Buster-report                  | : | wwPDB partial adaption of 1.1.7 (2018)                           |
| Percentile statistics          | : | 20250101.v01 (using entries in the PDB archive January 1st 2025) |
| Ideal geometry (proteins)      | : | Engh & Huber (2001)                                              |
| Ideal geometry (DNA, RNA)      | : | Parkinson et al. (1996)                                          |
| Validation Pipeline (wwPDB-VP) | : | 2.49                                                             |

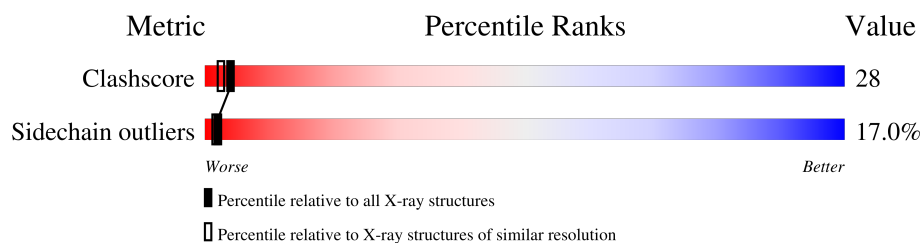
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 Å.

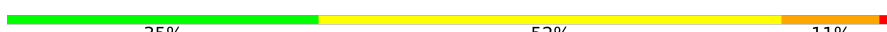
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



| Metric             | Whole archive<br>(#Entries) | Similar resolution<br>(#Entries, resolution range(Å)) |
|--------------------|-----------------------------|-------------------------------------------------------|
| Clashscore         | 190562                      | 1663 (2.36-2.36)                                      |
| Sidechain outliers | 187428                      | 1646 (2.36-2.36)                                      |

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ .

Note EDS was not executed.

| Mol | Chain | Length | Quality of chain                                                                     |
|-----|-------|--------|--------------------------------------------------------------------------------------|
| 1   | A     | 446    |  |
| 2   | B     | 439    |  |
| 3   | C     | 379    |  |
| 4   | D     | 241    |  |
| 5   | E     | 196    |  |
| 6   | F     | 110    |  |
| 7   | G     | 81     |  |
| 8   | H     | 78     |  |
| 9   | I     | 78     |  |

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| Mol | Chain | Length | Quality of chain                                                                   |
|-----|-------|--------|------------------------------------------------------------------------------------|
| 10  | J     | 62     |  |
| 11  | K     | 56     |  |

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

| Mol | Type | Chain | Res | Chirality | Geometry | Clashes | Electron density |
|-----|------|-------|-----|-----------|----------|---------|------------------|
| 14  | FES  | E     | 197 | -         | -        | X       | -                |

## 2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 16795 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called UBIQUINOL-CYTOCHROME C REDUCTASE COMPLEX CORE PROTEIN I.

| Mol | Chain | Residues | Atoms |      |     |     |    | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|---------|-------|
| 1   | A     | 446      | Total | C    | N   | O   | S  | 0       | 0       | 0     |
|     |       |          | 3458  | 2161 | 609 | 668 | 20 |         |         |       |

- Molecule 2 is a protein called UBIQUINOL-CYTOCHROME C REDUCTASE COMPLEX CORE PROTEIN 2.

| Mol | Chain | Residues | Atoms |      |     |     |   | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|---------|-------|
| 2   | B     | 423      | Total | C    | N   | O   | S | 0       | 0       | 0     |
|     |       |          | 3172  | 1993 | 562 | 610 | 7 |         |         |       |

- Molecule 3 is a protein called Cytochrome B.

| Mol | Chain | Residues | Atoms |      |     |     |    | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|---------|-------|
| 3   | C     | 377      | Total | C    | N   | O   | S  | 0       | 0       | 0     |
|     |       |          | 2995  | 2009 | 470 | 498 | 18 |         |         |       |

- Molecule 4 is a protein called Cytochrome c1, heme protein.

| Mol | Chain | Residues | Atoms |      |     |     |    | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|---------|-------|
| 4   | D     | 240      | Total | C    | N   | O   | S  | 0       | 0       | 0     |
|     |       |          | 1909  | 1219 | 328 | 347 | 15 |         |         |       |

- Molecule 5 is a protein called UBIQUINOL-CYTOCHROME C REDUCTASE IRON-SULFUR SUBUNIT.

| Mol | Chain | Residues | Atoms |     |     |     |   | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|---------|-------|
| 5   | E     | 196      | Total | C   | N   | O   | S | 0       | 0       | 0     |
|     |       |          | 1519  | 957 | 263 | 291 | 8 |         |         |       |

- Molecule 6 is a protein called Ubiquinol-cytochrome C reductase complex 14 kDa protein.

| Mol | Chain | Residues | Atoms |     |     |     |   | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|---------|-------|
| 6   | F     | 109      | Total | C   | N   | O   | S | 0       | 0       | 0     |
|     |       |          | 938   | 592 | 172 | 172 | 2 |         |         |       |

- Molecule 7 is a protein called Ubiquinol-cytochrome C reductase complex ubiquinone-binding protein QP-C.

| Mol | Chain | Residues | Atoms |     |     |     |   | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|---------|-------|
| 7   | G     | 79       | Total | C   | N   | O   | S | 0       | 0       | 0     |
|     |       |          | 662   | 431 | 123 | 107 | 1 |         |         |       |

- Molecule 8 is a protein called Ubiquinol-cytochrome C reductase complex 11 kDa protein.

| Mol | Chain | Residues | Atoms |     |     |     |   | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|---------|-------|
| 8   | H     | 78       | Total | C   | N   | O   | S | 0       | 0       | 0     |
|     |       |          | 639   | 384 | 111 | 139 | 5 |         |         |       |

- Molecule 9 is a protein called UBIQUINOL-CYTOCHROME C REDUCTASE 8 KDA PROTEIN.

| Mol | Chain | Residues | Atoms |     |    |    |   | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|---------|-------|
| 9   | I     | 57       | Total | C   | N  | O  | S | 0       | 0       | 0     |
|     |       |          | 406   | 253 | 77 | 74 | 2 |         |         |       |

- Molecule 10 is a protein called Ubiquinol-cytochrome C reductase complex 7.2 kDa protein.

| Mol | Chain | Residues | Atoms |     |    |    | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---------|---------|-------|
| 10  | J     | 60       | Total | C   | N  | O  | 0       | 0       | 0     |
|     |       |          | 495   | 324 | 86 | 85 |         |         |       |

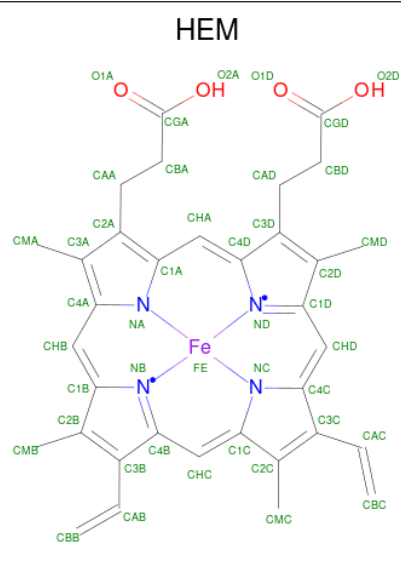
- Molecule 11 is a protein called Ubiquinol-cytochrome C reductase complex 6.4 kDa protein.

| Mol | Chain | Residues | Atoms |     |    |    |   | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|---------|-------|
| 11  | K     | 53       | Total | C   | N  | O  | S | 0       | 0       | 0     |
|     |       |          | 441   | 295 | 79 | 66 | 1 |         |         |       |

There is a discrepancy between the modelled and reference sequences:

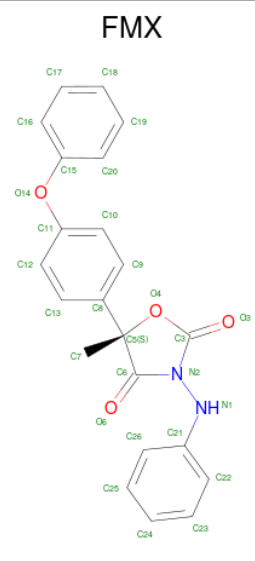
| Chain | Residue | Modelled | Actual | Comment  | Reference  |
|-------|---------|----------|--------|----------|------------|
| K     | 22      | GLN      | SER    | conflict | UNP P07552 |

- Molecule 12 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C<sub>34</sub>H<sub>32</sub>FeN<sub>4</sub>O<sub>4</sub>).



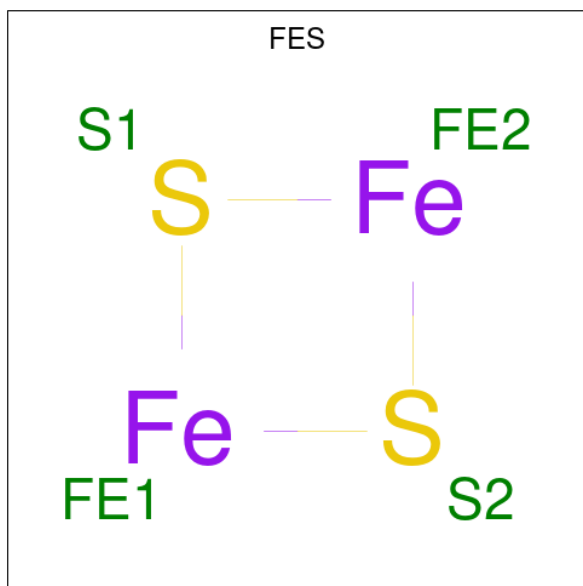
| Mol | Chain | Residues | Atoms       |         |         |        |        | ZeroOcc | AltConf |
|-----|-------|----------|-------------|---------|---------|--------|--------|---------|---------|
| 12  | C     | 1        | Total<br>43 | C<br>34 | Fe<br>1 | N<br>4 | O<br>4 | 0       | 0       |
| 12  | C     | 1        | Total<br>43 | C<br>34 | Fe<br>1 | N<br>4 | O<br>4 | 0       | 0       |
| 12  | D     | 1        | Total<br>43 | C<br>34 | Fe<br>1 | N<br>4 | O<br>4 | 0       | 0       |

- Molecule 13 is FAMOXADONE (CCD ID: FMX) (formula:  $C_{22}H_{18}N_2O_4$ ).



| Mol | Chain | Residues | Atoms |    |   |   | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---|---|---------|---------|
| 13  | C     | 1        | Total | C  | N | O | 0       | 0       |
|     |       |          | 28    | 22 | 2 | 4 |         |         |

- Molecule 14 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula:  $\text{Fe}_2\text{S}_2$ ).



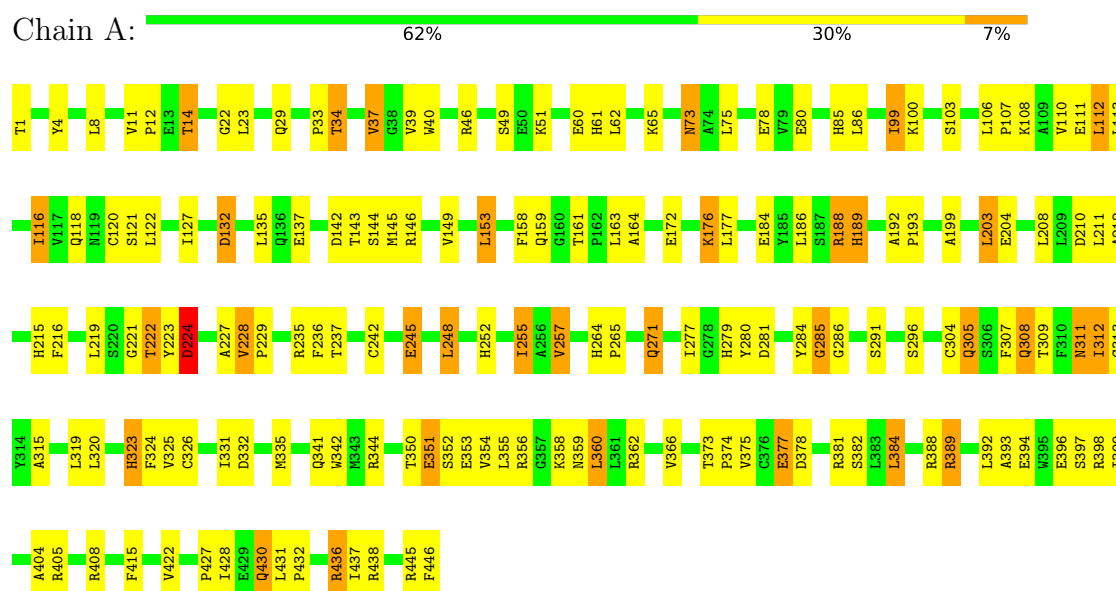
| Mol | Chain | Residues | Atoms |    |   | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---|---------|---------|
| 14  | E     | 1        | Total | Fe | S | 0       | 0       |
|     |       |          | 4     | 2  | 2 |         |         |

### 3 Residue-property plots

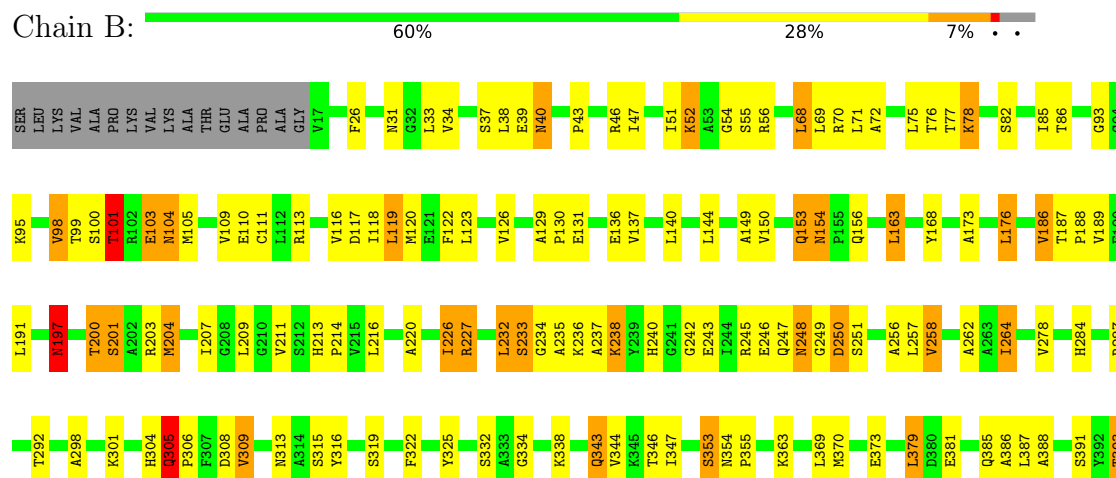
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

- Molecule 1: UBIQUINOL-CYTOCHROME C REDUCTASE COMPLEX CORE PROTEIN I



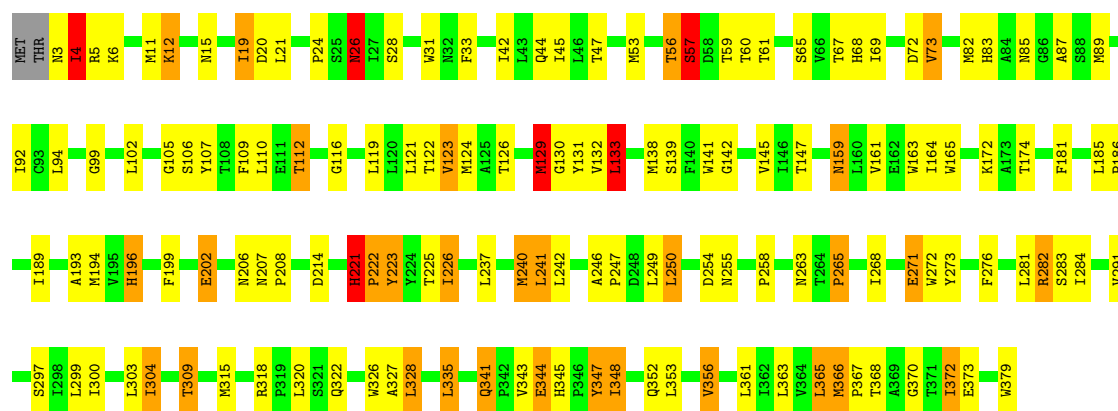
- Molecule 2: UBIQUINOL-CYTOCHROME C REDUCTASE COMPLEX CORE PROTEIN 2





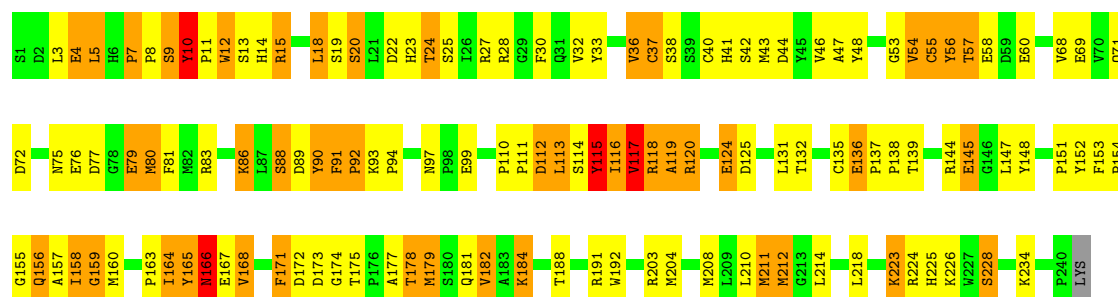
• Molecule 3: Cytochrome B

Chain C: 61% 29% 8% ..



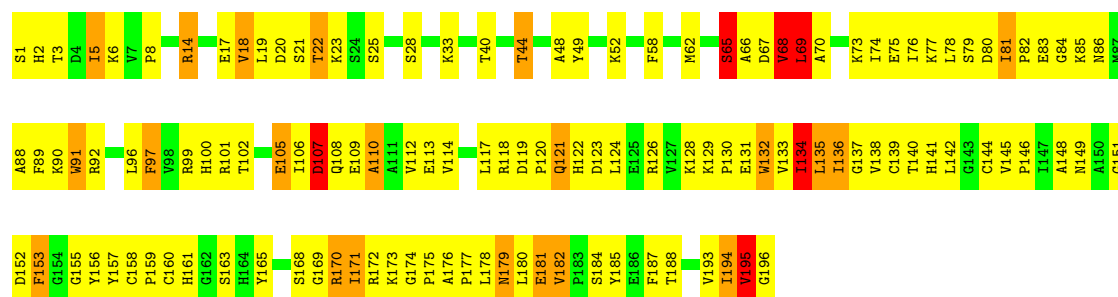
• Molecule 4: Cytochrome c1, heme protein

Chain D: 45% 34% 19% .



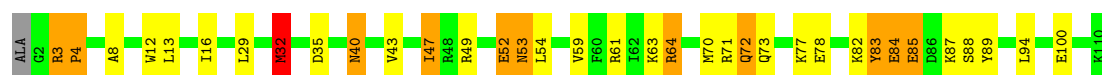
• Molecule 5: UBIQUINOL-CYTOCHROME C REDUCTASE IRON-SULFUR SUBUNIT

Chain E: 35% 52% 11% .



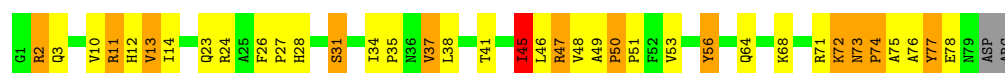
• Molecule 6: Ubiquinol-cytochrome C reductase complex 14 kDa protein

Chain F: 67% 21% 10% ..



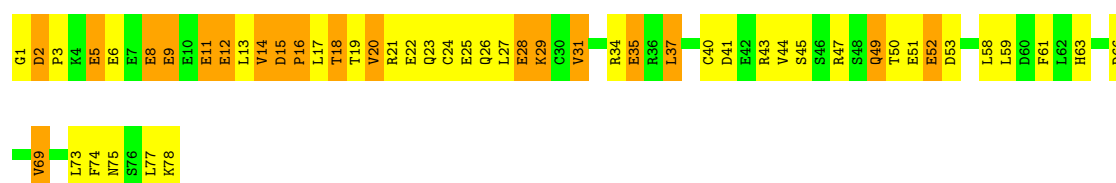
- Molecule 7: Ubiquinol-cytochrome C reductase complex ubiquinone-binding protein QP-C

Chain G: 52% 30% 15% ..



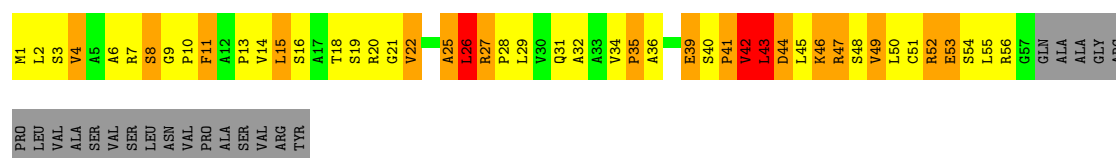
- Molecule 8: Ubiquinol-cytochrome C reductase complex 11 kDa protein

Chain H: 33% 42% 24%



- Molecule 9: UBIQUINOL-CYTOCHROME C REDUCTASE 8 KDA PROTEIN

Chain I: 13% 36% 21% • 27%



- Molecule 10: Ubiquinol-cytochrome C reductase complex 7.2 kDa protein

Chain J: 45% 37% 13% ..



- Molecule 11: Ubiquinol-cytochrome C reductase complex 6.4 kDa protein

Chain K: 61% 21% 12% 5%



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

| Property                                                 | Value                                           | Source    |
|----------------------------------------------------------|-------------------------------------------------|-----------|
| Space group                                              | I 41 2 2                                        | Depositor |
| Cell constants<br>a, b, c, $\alpha$ , $\beta$ , $\gamma$ | 154.09Å 154.09Å 591.84Å<br>90.00° 90.00° 90.00° | Depositor |
| Resolution (Å)                                           | 40.00 – 2.35                                    | Depositor |
| % Data completeness<br>(in resolution range)             | 96.0 (40.00-2.35)                               | Depositor |
| $R_{merge}$                                              | (Not available)                                 | Depositor |
| $R_{sym}$                                                | (Not available)                                 | Depositor |
| Refinement program                                       | REFMAC 5.0                                      | Depositor |
| R, $R_{free}$                                            | 0.259 , 0.306                                   | Depositor |
| Estimated twinning fraction                              | No twinning to report.                          | Xtriage   |
| Total number of atoms                                    | 16795                                           | wwPDB-VP  |
| Average B, all atoms (Å <sup>2</sup> )                   | 25.0                                            | wwPDB-VP  |

## 5 Model quality (i)

### 5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, FMX, FES

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

| Mol | Chain | Bond lengths |                 | Bond angles |                  |
|-----|-------|--------------|-----------------|-------------|------------------|
|     |       | RMSZ         | # $ Z  > 5$     | RMSZ        | # $ Z  > 5$      |
| 1   | A     | 1.11         | 3/3531 (0.1%)   | 1.14        | 16/4792 (0.3%)   |
| 2   | B     | 1.44         | 16/3232 (0.5%)  | 1.29        | 23/4386 (0.5%)   |
| 3   | C     | 1.11         | 5/3092 (0.2%)   | 1.24        | 33/4232 (0.8%)   |
| 4   | D     | 0.96         | 4/1968 (0.2%)   | 1.26        | 23/2673 (0.9%)   |
| 5   | E     | 1.01         | 2/1553 (0.1%)   | 1.37        | 21/2100 (1.0%)   |
| 6   | F     | 1.27         | 3/958 (0.3%)    | 1.19        | 7/1284 (0.5%)    |
| 7   | G     | 1.12         | 0/684           | 1.23        | 7/926 (0.8%)     |
| 8   | H     | 0.84         | 0/645           | 1.41        | 7/864 (0.8%)     |
| 9   | I     | 1.39         | 5/411 (1.2%)    | 1.85        | 12/558 (2.2%)    |
| 10  | J     | 0.90         | 1/508 (0.2%)    | 1.19        | 5/686 (0.7%)     |
| 11  | K     | 0.99         | 0/457           | 1.14        | 1/625 (0.2%)     |
| All | All   | 1.16         | 39/17039 (0.2%) | 1.26        | 155/23126 (0.7%) |

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

| Mol | Chain | #Chirality outliers | #Planarity outliers |
|-----|-------|---------------------|---------------------|
| 3   | C     | 0                   | 2                   |

The worst 5 of 39 bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z      | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|-------|--------|-------------|----------|
| 2   | B     | 424 | MET  | SD-CE | -16.77 | 1.37        | 1.79     |
| 4   | D     | 211 | MET  | SD-CE | -10.21 | 1.54        | 1.79     |
| 3   | C     | 129 | MET  | SD-CE | -9.66  | 1.55        | 1.79     |
| 9   | I     | 43  | LEU  | N-CA  | -8.34  | 1.35        | 1.46     |
| 4   | D     | 119 | ALA  | CA-C  | -7.92  | 1.42        | 1.52     |

The worst 5 of 155 bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms  | Z      | Observed(°) | Ideal(°) |
|-----|-------|-----|------|--------|--------|-------------|----------|
| 2   | B     | 233 | SER  | N-CA-C | -14.19 | 92.88       | 110.61   |
| 9   | I     | 42  | VAL  | O-C-N  | 14.04  | 140.12      | 122.57   |
| 3   | C     | 221 | HIS  | CA-C-N | 12.36  | 135.29      | 119.84   |
| 3   | C     | 221 | HIS  | C-N-CA | 12.36  | 135.29      | 119.84   |
| 4   | D     | 118 | ARG  | N-CA-C | 12.09  | 125.85      | 111.82   |

There are no chirality outliers.

All (2) planarity outliers are listed below:

| Mol | Chain | Res | Type | Group             |
|-----|-------|-----|------|-------------------|
| 3   | C     | 221 | HIS  | Mainchain,Peptide |

## 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 1   | A     | 3458  | 0        | 3356     | 152     | 0            |
| 2   | B     | 3172  | 0        | 3152     | 123     | 0            |
| 3   | C     | 2995  | 0        | 3058     | 125     | 0            |
| 4   | D     | 1909  | 0        | 1857     | 175     | 0            |
| 5   | E     | 1519  | 0        | 1503     | 182     | 0            |
| 6   | F     | 938   | 0        | 932      | 28      | 0            |
| 7   | G     | 662   | 0        | 662      | 34      | 0            |
| 8   | H     | 639   | 0        | 604      | 50      | 0            |
| 9   | I     | 406   | 0        | 437      | 99      | 0            |
| 10  | J     | 495   | 0        | 493      | 69      | 0            |
| 11  | K     | 441   | 0        | 450      | 27      | 0            |
| 12  | C     | 86    | 0        | 60       | 22      | 0            |
| 12  | D     | 43    | 0        | 30       | 6       | 0            |
| 13  | C     | 28    | 0        | 18       | 0       | 0            |
| 14  | E     | 4     | 0        | 0        | 2       | 0            |
| All | All   | 16795 | 0        | 16612    | 933     | 0            |

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 28.

The worst 5 of 933 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

| Atom-1         | Atom-2           | Interatomic distance (Å) | Clash overlap (Å) |
|----------------|------------------|--------------------------|-------------------|
| 10:J:49:GLY:C  | 10:J:54:HIS:HE1  | 1.40                     | 1.28              |
| 4:D:40:CYS:SG  | 12:D:242:HEM:HAC | 1.76                     | 1.26              |
| 10:J:51:LEU:O  | 10:J:55:ILE:HG13 | 1.43                     | 1.18              |
| 3:C:129:MET:CE | 3:C:181:PHE:HB2  | 1.75                     | 1.17              |
| 9:I:47:ARG:HG2 | 9:I:48:SER:H     | 1.09                     | 1.16              |

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

| Mol | Chain | Analysed        | Rotameric  | Outliers  | Percentiles |   |
|-----|-------|-----------------|------------|-----------|-------------|---|
| 1   | A     | 370/370 (100%)  | 315 (85%)  | 55 (15%)  | 3           | 2 |
| 2   | B     | 332/343 (97%)   | 288 (87%)  | 44 (13%)  | 4           | 3 |
| 3   | C     | 325/327 (99%)   | 279 (86%)  | 46 (14%)  | 3           | 3 |
| 4   | D     | 205/206 (100%)  | 159 (78%)  | 46 (22%)  | 1           | 1 |
| 5   | E     | 168/168 (100%)  | 139 (83%)  | 29 (17%)  | 2           | 1 |
| 6   | F     | 98/98 (100%)    | 80 (82%)   | 18 (18%)  | 1           | 1 |
| 7   | G     | 69/71 (97%)     | 53 (77%)   | 16 (23%)  | 1           | 0 |
| 8   | H     | 74/74 (100%)    | 56 (76%)   | 18 (24%)  | 1           | 0 |
| 9   | I     | 44/60 (73%)     | 30 (68%)   | 14 (32%)  | 0           | 0 |
| 10  | J     | 50/52 (96%)     | 42 (84%)   | 8 (16%)   | 2           | 2 |
| 11  | K     | 43/46 (94%)     | 35 (81%)   | 8 (19%)   | 1           | 1 |
| All | All   | 1778/1815 (98%) | 1476 (83%) | 302 (17%) | 2           | 2 |

5 of 302 residues with a non-rotameric sidechain are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 6   | F     | 77  | LYS  |
| 10  | J     | 5   | LEU  |
| 7   | G     | 2   | ARG  |
| 8   | H     | 18  | THR  |
| 11  | K     | 34  | TRP  |

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 73 such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 5   | E     | 116 | GLN  |
| 11  | K     | 12  | GLN  |
| 5   | E     | 161 | HIS  |
| 6   | F     | 73  | GLN  |
| 2   | B     | 162 | ASN  |

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

| Mol | Type | Chain | Res | Link | Bond lengths |      |          | Bond angles |      |          |
|-----|------|-------|-----|------|--------------|------|----------|-------------|------|----------|
|     |      |       |     |      | Counts       | RMSZ | # Z  > 2 | Counts      | RMSZ | # Z  > 2 |
| 14  | FES  | E     | 197 | 5    | 0,4,4        | -    | -        | -           |      |          |
| 12  | HEM  | D     | 242 | 4    | 50,50,50     | 1.82 | 7 (14%)  | 67,82,82    | 1.46 | 11 (16%) |
| 12  | HEM  | C     | 610 | 3    | 50,50,50     | 3.40 | 28 (56%) | 67,82,82    | 2.67 | 30 (44%) |
| 12  | HEM  | C     | 609 | 3    | 50,50,50     | 3.56 | 28 (56%) | 67,82,82    | 2.63 | 32 (47%) |
| 13  | FMX  | C     | 611 | -    | 30,31,31     | 0.94 | 2 (6%)   | 36,44,44    | 1.23 | 4 (11%)  |

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

| Mol | Type | Chain | Res | Link | Chirals | Torsions   | Rings   |
|-----|------|-------|-----|------|---------|------------|---------|
| 14  | FES  | E     | 197 | 5    | -       | -          | 0/1/1/1 |
| 12  | HEM  | D     | 242 | 4    | -       | 6/14/54/54 | -       |
| 12  | HEM  | C     | 610 | 3    | -       | 7/14/54/54 | -       |
| 12  | HEM  | C     | 609 | 3    | -       | 9/14/54/54 | -       |
| 13  | FMX  | C     | 611 | -    | -       | 2/14/33/33 | 0/4/4/4 |

The worst 5 of 65 bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms   | Z     | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|---------|-------|-------------|----------|
| 12  | C     | 610 | HEM  | C3C-C4C | -7.50 | 1.32        | 1.46     |
| 12  | C     | 609 | HEM  | C3B-C4B | -7.22 | 1.30        | 1.44     |
| 12  | D     | 242 | HEM  | C3D-C2D | 7.14  | 1.52        | 1.36     |
| 12  | C     | 609 | HEM  | CHA-C4D | 6.80  | 1.52        | 1.38     |
| 12  | C     | 609 | HEM  | CHD-C4C | 6.54  | 1.51        | 1.38     |

The worst 5 of 77 bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms       | Z     | Observed(°) | Ideal(°) |
|-----|-------|-----|------|-------------|-------|-------------|----------|
| 12  | C     | 610 | HEM  | C4C-CHD-C1D | -6.73 | 111.72      | 126.02   |
| 12  | C     | 609 | HEM  | C4C-CHD-C1D | -6.33 | 112.57      | 126.02   |
| 12  | C     | 609 | HEM  | C1C-CHC-C4B | -6.08 | 113.09      | 126.02   |
| 12  | C     | 610 | HEM  | C1C-CHC-C4B | -6.05 | 113.15      | 126.02   |
| 12  | C     | 609 | HEM  | CHB-C4A-NA  | 5.16  | 133.22      | 123.86   |

There are no chirality outliers.

5 of 24 torsion outliers are listed below:

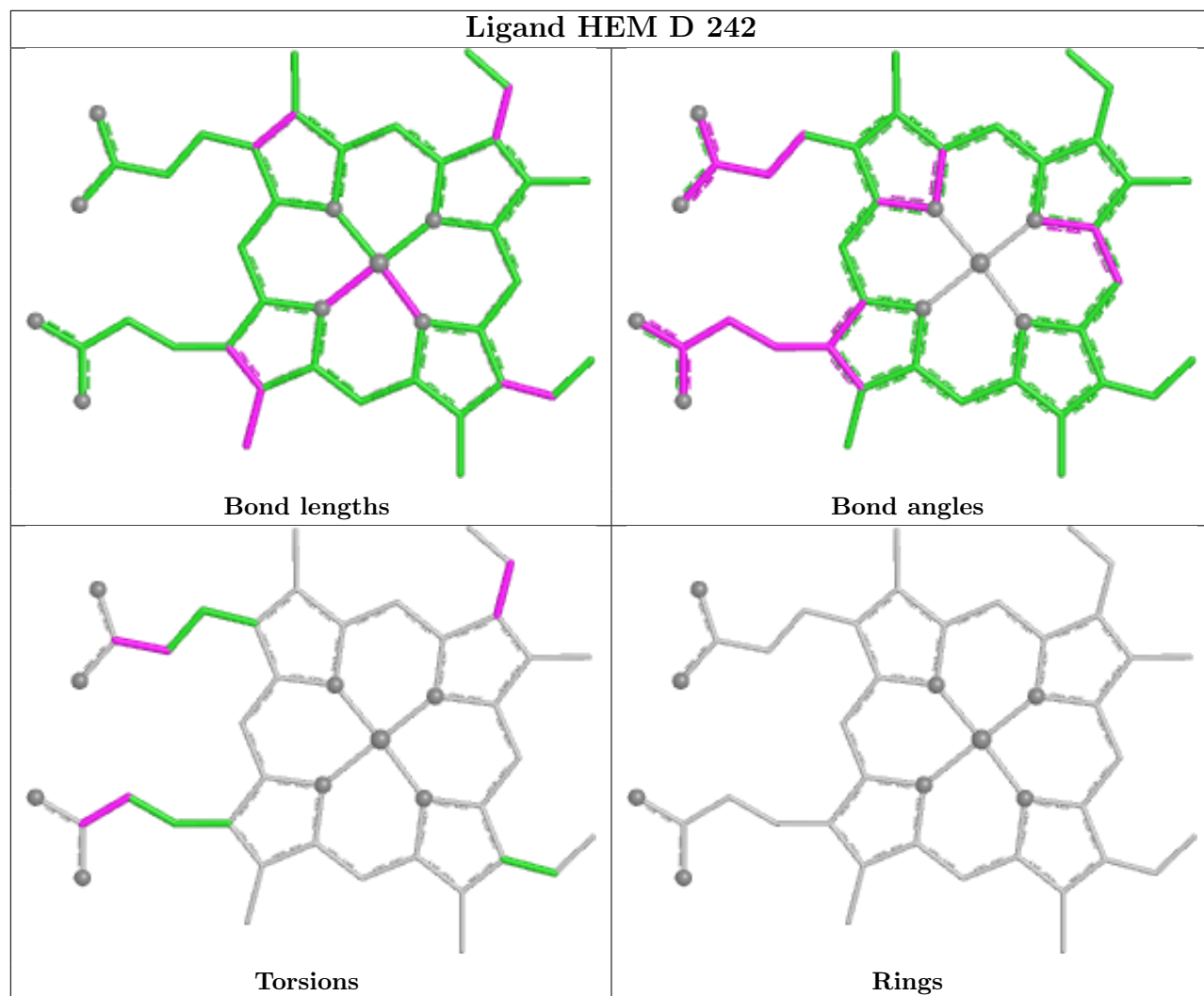
| Mol | Chain | Res | Type | Atoms           |
|-----|-------|-----|------|-----------------|
| 12  | C     | 609 | HEM  | C2D-C3D-CAD-CBD |
| 12  | C     | 609 | HEM  | C4D-C3D-CAD-CBD |
| 12  | C     | 610 | HEM  | C2B-C3B-CAB-CBB |
| 12  | C     | 610 | HEM  | C2C-C3C-CAC-CBC |
| 12  | C     | 610 | HEM  | C4C-C3C-CAC-CBC |

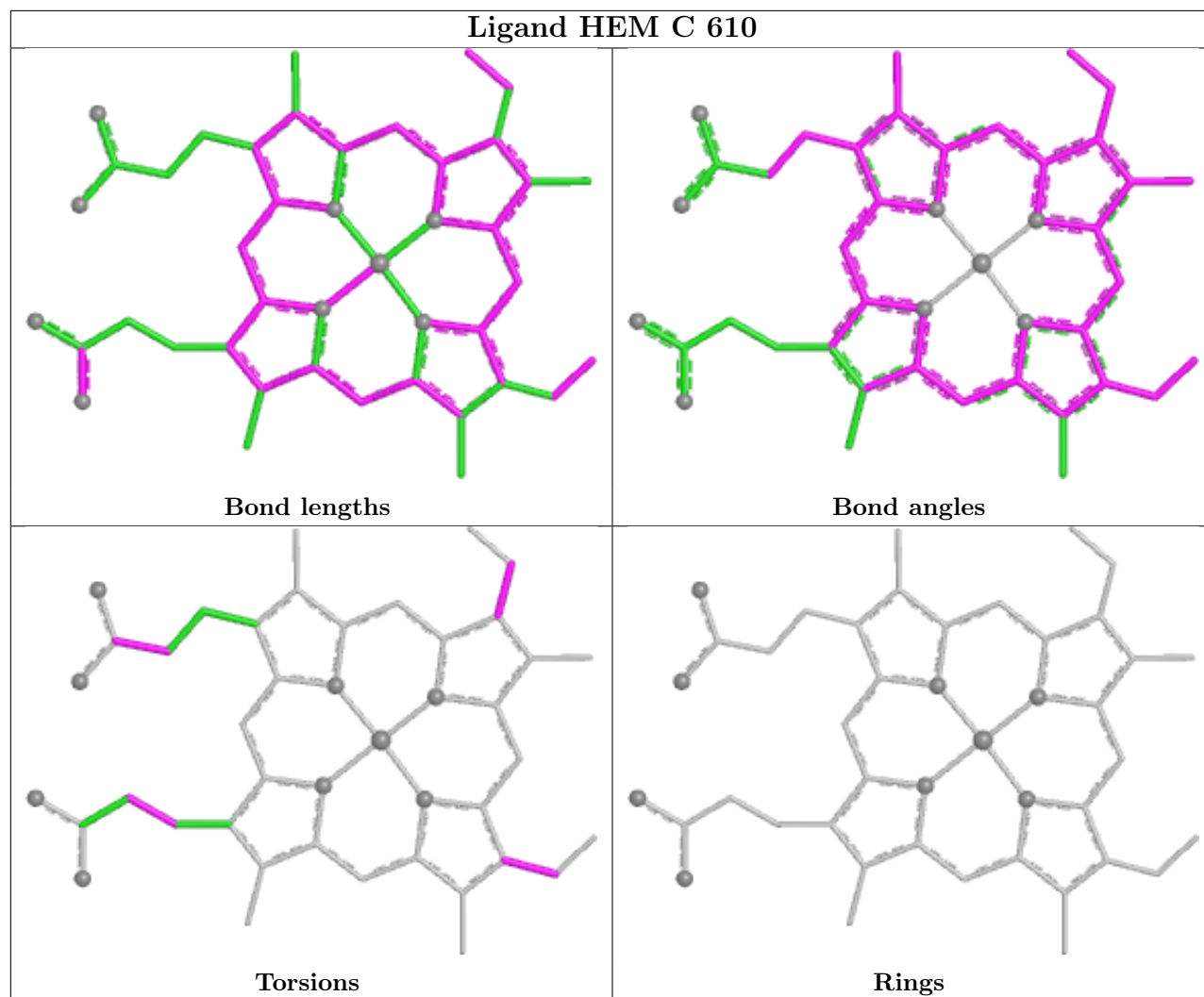
There are no ring outliers.

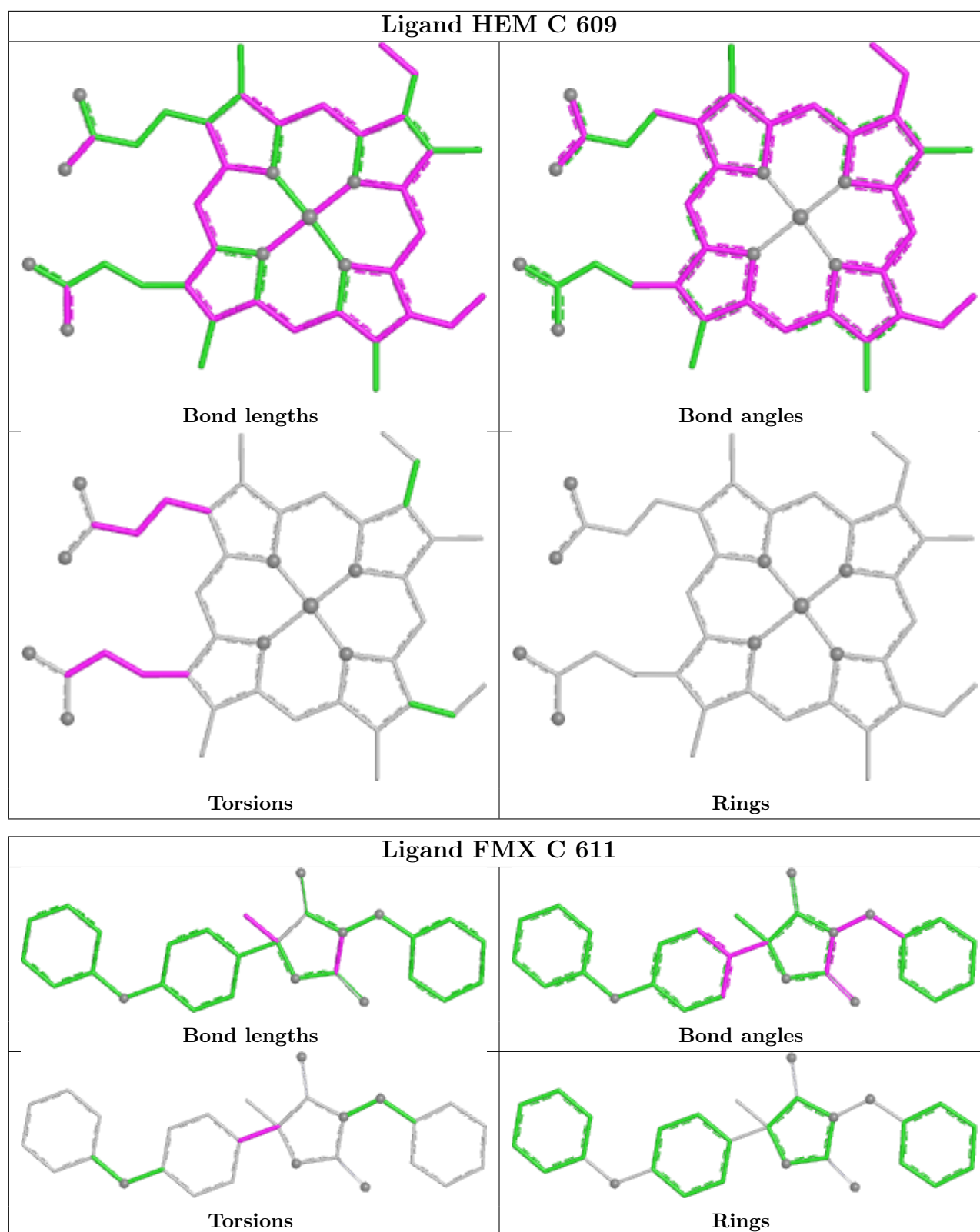
4 monomers are involved in 30 short contacts:

| Mol | Chain | Res | Type | Clashes | Symm-Clashes |
|-----|-------|-----|------|---------|--------------|
| 14  | E     | 197 | FES  | 2       | 0            |
| 12  | D     | 242 | HEM  | 6       | 0            |
| 12  | C     | 610 | HEM  | 11      | 0            |
| 12  | C     | 609 | HEM  | 11      | 0            |

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.