



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 8, 2025 – 05:47 pm BST

PDB ID : 9HQ1 / pdb_00009hq1
Title : XusB lipoprotein bound to ferric salmochelin
Authors : Silale, A.; Soo, Y.L.; van den Berg, B.
Deposited on : 2024-12-16
Resolution : 1.80 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

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>.

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

MolProbity : **FAILED**
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

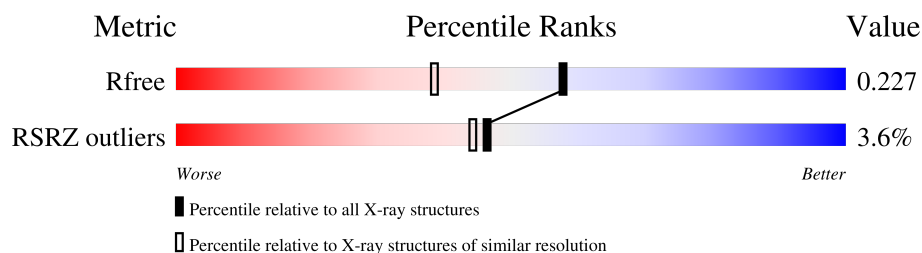
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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 (#Entries) | Similar resolution (#Entries, resolution range(Å)) |
|---------------|-----------------------------|---|
| R_{free} | 164625 | 7108 (1.80-1.80) |
| RSRZ outliers | 164620 | 7108 (1.80-1.80) |

MolProbity failed to run properly - the sequence quality summary graphics cannot be shown.

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7471 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.

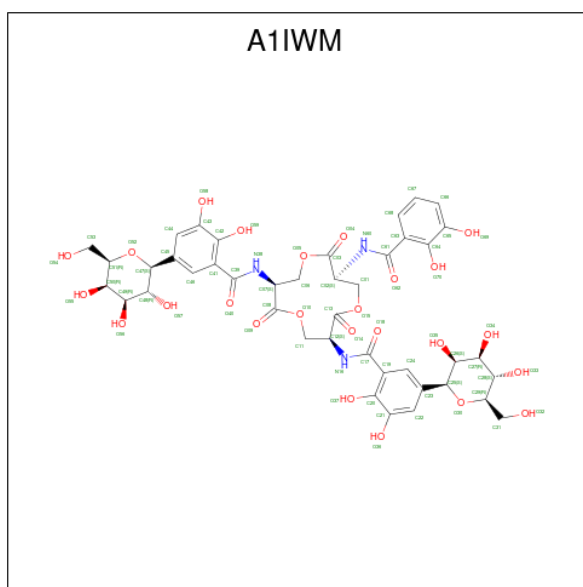
- Molecule 1 is a protein called DUF4374 domain-containing protein.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|---------|-------|
| 1 | A | 426 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 3307 | 2090 | 547 | 660 | 10 | | | |
| 1 | B | 426 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 3298 | 2084 | 545 | 659 | 10 | | | |

There are 16 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------------|------------|
| A | 465 | LEU | - | expression tag | UNP Q8A622 |
| A | 466 | GLU | - | expression tag | UNP Q8A622 |
| A | 467 | HIS | - | expression tag | UNP Q8A622 |
| A | 468 | HIS | - | expression tag | UNP Q8A622 |
| A | 469 | HIS | - | expression tag | UNP Q8A622 |
| A | 470 | HIS | - | expression tag | UNP Q8A622 |
| A | 471 | HIS | - | expression tag | UNP Q8A622 |
| A | 472 | HIS | - | expression tag | UNP Q8A622 |
| B | 465 | LEU | - | expression tag | UNP Q8A622 |
| B | 466 | GLU | - | expression tag | UNP Q8A622 |
| B | 467 | HIS | - | expression tag | UNP Q8A622 |
| B | 468 | HIS | - | expression tag | UNP Q8A622 |
| B | 469 | HIS | - | expression tag | UNP Q8A622 |
| B | 470 | HIS | - | expression tag | UNP Q8A622 |
| B | 471 | HIS | - | expression tag | UNP Q8A622 |
| B | 472 | HIS | - | expression tag | UNP Q8A622 |

- Molecule 2 is Salmochelin S4 (CCD ID: A1IWM) (formula: C₄₂H₄₇N₃O₂₅) (labeled as "Ligand of Interest" by depositor).



| Mol | Chain | Residues | Atoms | | | | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---|----|---------|---------|
| 2 | A | 1 | Total | C | N | O | 0 | 0 |
| | | | 70 | 42 | 3 | 25 | | |

- Molecule 3 is FE (III) ION (CCD ID: FE) (formula: Fe).

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---------|---------|
| 3 | A | 1 | Total | Fe | 0 | 0 |
| | | | 1 | 1 | | |

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---------|---------|
| 4 | A | 2 | Total | Ca | 0 | 0 |
| | | | 2 | 2 | | |
| 4 | B | 4 | Total | Ca | 0 | 0 |
| | | | 4 | 4 | | |

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---------|---------|
| 5 | A | 2 | Total | Cl | 0 | 0 |
| | | | 2 | 2 | | |
| 5 | B | 2 | Total | Cl | 0 | 0 |
| | | | 2 | 2 | | |

- Molecule 6 is water.

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|--------------|----------|---------|---------|
| 6 | A | 384 | Total 384 | O 384 | 0 | 0 |
| 6 | B | 401 | Total 401 | O 401 | 0 | 0 |

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3 Data and refinement statistics

| Property | Value | Source |
|---|---|------------------|
| Space group | P 1 21 1 | Depositor |
| Cell constants a, b, c, α , β , γ | 45.44Å 108.72Å 93.83Å 90.00° 96.66° 90.00° | Depositor |
| Resolution (Å) | 54.36 – 1.80 54.36 – 1.80 | Depositor EDS |
| % Data completeness (in resolution range) | 99.8 (54.36-1.80) 99.7 (54.36-1.80) | Depositor EDS |
| R_{merge} | 0.10 | Depositor |
| R_{sym} | (Not available) | Depositor |
| $\langle I/\sigma(I) \rangle$ ¹ | 1.66 (at 1.80Å) | Xtriage |
| Refinement program | PHENIX 1.21.1_5286 | Depositor |
| R, R_{free} | 0.193 , 0.229 0.193 , 0.227 | Depositor DCC |
| R_{free} test set | 4149 reflections (4.96%) | wwPDB-VP |
| Wilson B-factor (Å ²) | 23.3 | Xtriage |
| Anisotropy | 0.213 | Xtriage |
| Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²) | 0.34 , 32.9 | EDS |
| L-test for twinning ² | $\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$ | Xtriage |
| Estimated twinning fraction | No twinning to report. | Xtriage |
| F_o, F_c correlation | 0.95 | EDS |
| Total number of atoms | 7471 | wwPDB-VP |
| Average B, all atoms (Å ²) | 26.0 | wwPDB-VP |

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

4 Model quality [i](#)

4.1 Standard geometry [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.2 Too-close contacts [i](#)

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4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

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4.3.2 Protein sidechains [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3.3 RNA [i](#)

MolProbity failed to run properly - this section is therefore empty.

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

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

4.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

4.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 11 are monoatomic - leaving 1 for Mogul analysis.

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 | $\# Z > 2$ |
| 2 | A1IWM | A | 501 | 3 | 75,75,75 | 2.07 | 22 (29%) | 110,110,110 | 1.92 | 32 (29%) |

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 |
|-----|-------|-------|-----|------|---------|--------------|---------|
| 2 | A1IWM | A | 501 | 3 | - | 2/63/103/103 | 0/5/6/6 |

The worst 5 of 22 bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(Å) | Ideal(Å) |
|-----|-------|-----|-------|---------|-------|-------------|----------|
| 2 | A | 501 | A1IWM | C26-C25 | 5.31 | 1.65 | 1.53 |
| 2 | A | 501 | A1IWM | C17-N16 | 5.20 | 1.45 | 1.34 |
| 2 | A | 501 | A1IWM | O05-C03 | 4.31 | 1.42 | 1.33 |
| 2 | A | 501 | A1IWM | O52-C47 | -4.19 | 1.38 | 1.44 |
| 2 | A | 501 | A1IWM | C61-N60 | 4.08 | 1.43 | 1.34 |

The worst 5 of 32 bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | Ideal(°) |
|-----|-------|-----|-------|-------------|-------|-------------|----------|
| 2 | A | 501 | A1IWM | C23-C25-C26 | 6.22 | 127.81 | 112.62 |
| 2 | A | 501 | A1IWM | O52-C47-C45 | 5.74 | 117.96 | 108.00 |
| 2 | A | 501 | A1IWM | O30-C29-C31 | 5.00 | 118.86 | 106.44 |
| 2 | A | 501 | A1IWM | O30-C25-C23 | -4.18 | 100.74 | 108.00 |
| 2 | A | 501 | A1IWM | O10-C08-O09 | -4.14 | 116.28 | 124.13 |

There are no chirality outliers.

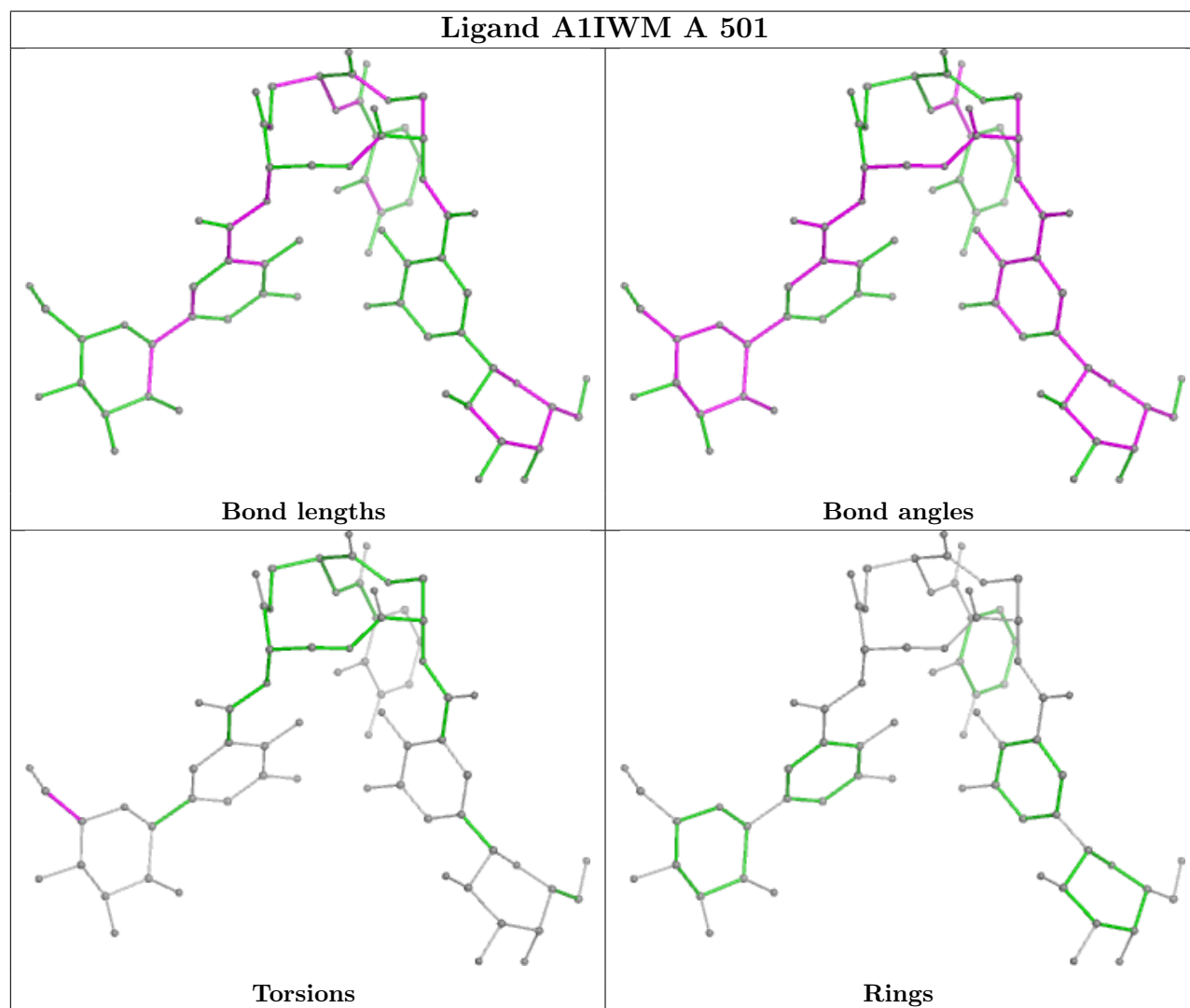
All (2) torsion outliers are listed below:

| Mol | Chain | Res | Type | Atoms |
|-----|-------|-----|-------|-----------------|
| 2 | A | 501 | A1IWM | C28-C29-C31-O32 |
| 2 | A | 501 | A1IWM | O30-C29-C31-O32 |

There are no ring outliers.

No monomer is involved in short contacts.

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.



4.7 Other polymers ⓘ

There are no such residues in this entry.

4.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

| Mol | Chain | Analysed | <RSRZ> | #RSRZ>2 | OWAB(Å ²) | Q<0.9 |
|-----|-------|---------------|--------|---------------|-----------------------|--------|
| 1 | A | 426/472 (90%) | -0.06 | 10 (2%) 61 59 | 11, 24, 40, 53 | 1 (0%) |
| 1 | B | 426/472 (90%) | -0.04 | 21 (4%) 36 33 | 16, 23, 44, 63 | 0 |
| All | All | 852/944 (90%) | -0.05 | 31 (3%) 46 44 | 11, 24, 42, 63 | 1 (0%) |

The worst 5 of 31 RSRZ outliers are listed below:

| Mol | Chain | Res | Type | RSRZ |
|-----|-------|-----|------|------|
| 1 | B | 418 | VAL | 5.9 |
| 1 | B | 417 | GLY | 4.0 |
| 1 | B | 415 | ASN | 3.6 |
| 1 | B | 411 | TYR | 3.2 |
| 1 | A | 388 | ASN | 3.2 |

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

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

5.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.4 Ligands [i](#)

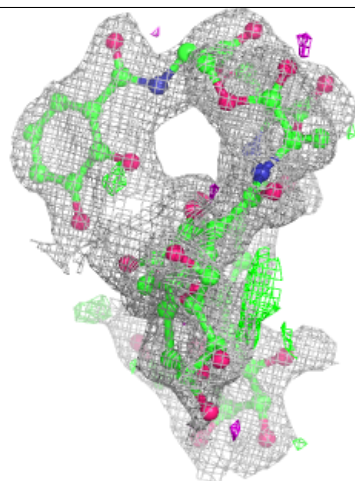
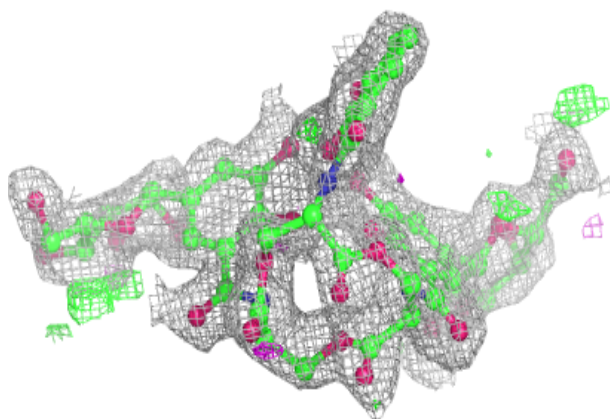
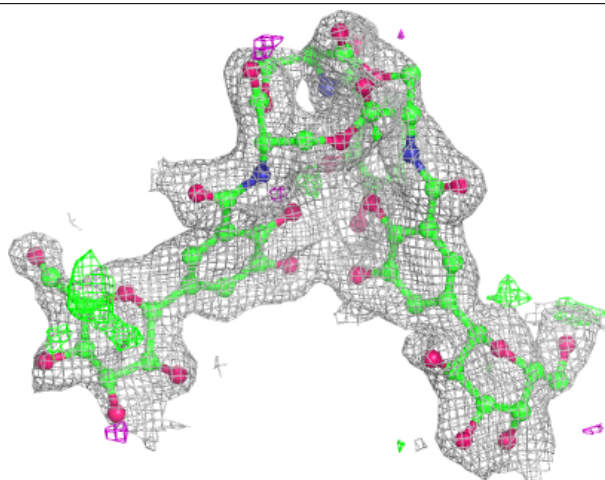
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

| Mol | Type | Chain | Res | Atoms | RSCC | RSR | B-factors(\AA^2) | Q<0.9 |
|-----|-------|-------|-----|-------|------|------|-----------------------------|-------|
| 2 | A1IWM | A | 501 | 70/70 | 0.86 | 0.11 | 26,35,40,42 | 70 |
| 5 | CL | A | 505 | 1/1 | 0.93 | 0.10 | 40,40,40,40 | 0 |
| 4 | CA | B | 502 | 1/1 | 0.95 | 0.11 | 42,42,42,42 | 0 |
| 4 | CA | B | 504 | 1/1 | 0.96 | 0.11 | 49,49,49,49 | 0 |
| 4 | CA | A | 504 | 1/1 | 0.97 | 0.12 | 43,43,43,43 | 0 |
| 5 | CL | B | 506 | 1/1 | 0.97 | 0.07 | 47,47,47,47 | 0 |
| 4 | CA | B | 503 | 1/1 | 0.98 | 0.16 | 41,41,41,41 | 0 |
| 4 | CA | B | 501 | 1/1 | 0.98 | 0.07 | 27,27,27,27 | 0 |
| 4 | CA | A | 503 | 1/1 | 0.99 | 0.03 | 24,24,24,24 | 0 |
| 5 | CL | A | 506 | 1/1 | 0.99 | 0.03 | 25,25,25,25 | 0 |
| 5 | CL | B | 505 | 1/1 | 0.99 | 0.04 | 26,26,26,26 | 0 |
| 3 | FE | A | 502 | 1/1 | 0.99 | 0.03 | 27,27,27,27 | 0 |

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around A1IWM A 501:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



5.5 Other polymers [i](#)

There are no such residues in this entry.