

wwPDB X-ray Structure Validation Summary Report (i)

Feb 21, 2024 – 09:03 AM EST

PDB ID : 4O35

Title: Crystal structure of carbomonoxy murine neuroglobin mutant F106W

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Deposited on : 2013-12-18

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 (i) 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 (1)) were used in the production of this report:

MolProbity: 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13

EDS : 2.36

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

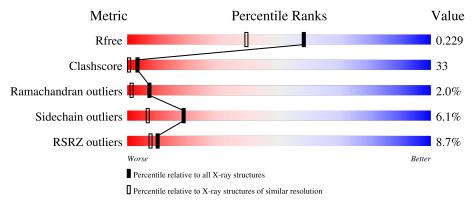
Validation Pipeline (wwPDB-VP) : 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\ DIFFRACTION$

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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\mathring{\rm A})}) \end{array}$
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 >=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 <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain		
		1 = 4	8%		
1	A	154	75%	16%	5% • •



2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 1574 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.

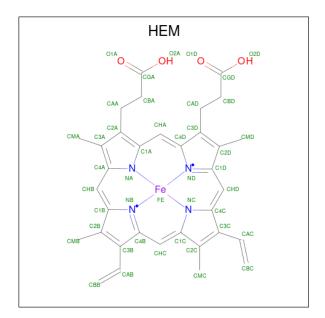
• Molecule 1 is a protein called Neuroglobin.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Δ	149	Total	С	N	О	S	0	36	0
1	11	143	1420	916	246	253	5		30	U

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q9ER97
A	-1	SER	-	expression tag	UNP Q9ER97
A	0	HIS	-	expression tag	UNP Q9ER97
A	55	SER	CYS	engineered mutation	UNP Q9ER97
A	106	TRP	PHE	engineered mutation	UNP Q9ER97
A	120	SER	CYS	engineered mutation	UNP Q9ER97

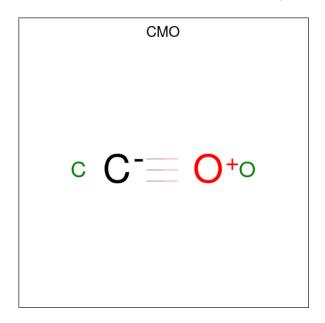
• Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).





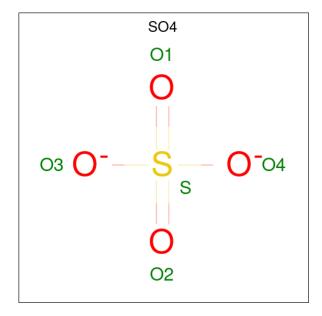
Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf
2	Λ	1	Total	С	Fe	N	О	0	0
	A	1	43	34	1	4	4	0	0

• Molecule 3 is CARBON MONOXIDE (three-letter code: CMO) (formula: CO).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
3	A	1	Total 2	C 1	O 1	0	0

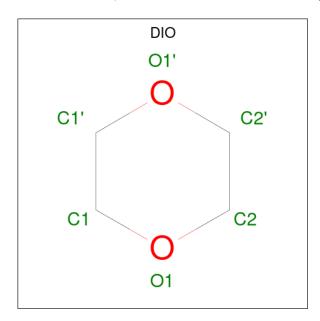
 \bullet Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: $\mathrm{O_4S}).$





\mathbf{Mol}	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0

 \bullet Molecule 5 is 1,4-DIETHYLENE DIOXIDE (three-letter code: DIO) (formula: $\mathrm{C_4H_8O_2}).$



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 6 4 2	0	0
5	A	1	Total C O 6 4 2	0	0
5	A	1	Total C O 6 4 2	0	0

• Molecule 6 is water.

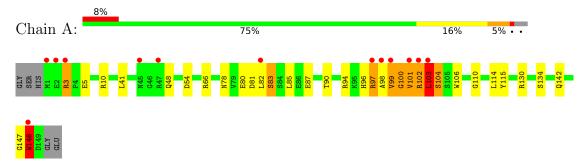
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	86	Total O 86 86	0	0



3 Residue-property plots (i)

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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). 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.

• Molecule 1: Neuroglobin





4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants	89.16Å 89.16Å 108.62Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.18 - 1.80	Depositor
Resolution (A)	28.18 - 1.80	EDS
% Data completeness	99.1 (28.18-1.80)	Depositor
(in resolution range)	99.1 (28.18-1.80)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	7.47 (at 1.80Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
P.P.	0.188 , 0.234	Depositor
R, R_{free}	0.194 , 0.229	DCC
R_{free} test set	1545 reflections $(9.98%)$	wwPDB-VP
Wilson B-factor (Å ²)	21.1	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34, 51.9	EDS
L-test for twinning ²	$< L > = 0.50, < L^2> = 0.33$	Xtriage
	0.029 for -2/3 +h- 1/3 +k+ 2/3 +l,- 1/3 +h- 2/3 +k-	
	2/3*1,2/3*h-2/3*k+1/3*l	
Estimated twinning fraction	0.022 for $-h,1/3*h-1/3*k+2/3*1,2/3*h+4/3*$	Xtriage
	k+1/3*l 0.015 for -1/3*h+1/3*k-2/3*l,-k,-4/3*h-2/3	
	*k+1/3*l	
F_o, F_c correlation	0.95	EDS
Total number of atoms	1574	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 8.54% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities estimated from amplitudes.

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: CMO, SO4, DIO, HEM

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).

Mal	Chain	Bond	lengths	Bond angles		
MIOI	Chain	RMSZ $ \# Z > 5$		RMSZ $\# Z > 5$		
1	A	0.51	0/1533	0.66	2/2078 (0.1%)	

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 maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\mathrm{Ideal}(^{o})$
1	A	100[A]	GLY	N-CA-C	5.09	125.83	113.10
1	A	100[B]	GLY	N-CA-C	5.09	125.83	113.10

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	102[B]	ARG	Peptide
1	A	103[B]	LEU	Peptide
1	A	148[B]	TRP	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	1420	0	1472	97	2
2	A	43	0	30	2	0
3	A	2	0	0	0	0
4	A	5	0	0	0	0
5	A	18	0	24	2	0
6	A	86	0	0	2	1
All	All	1574	0	1526	98	2

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 33.

The worst 5 of 98 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:97[A]:ARG:CB	1:A:98[A]:ALA:HB2	1.39	1.50
1:A:102[B]:ARG:CB	1:A:103[B]:LEU:HG	1.43	1.44
1:A:102[B]:ARG:HB3	1:A:103[B]:LEU:CG	1.56	1.33
1:A:98[A]:ALA:O	1:A:100[A]:GLY:N	1.73	1.21
1:A:102[B]:ARG:CA	1:A:103[B]:LEU:HG	1.69	1.21

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:A:54:ASP:OD1	1:A:115[C]:TYR:OH[5_555]	1.95	0.25
1:A:66[B]:ARG:NH1	6:A:372:HOH:O[5_555]	2.16	0.04

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran 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 backbone conformation was analysed, and the total number of residues.



Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	183/154 (119%)	170 (93%)	8 (4%)	5 (3%)	5 1

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	103[A]	LEU
1	A	103[B]	LEU
1	A	83	SER
1	A	101[A]	VAL
1	A	101[B]	VAL

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	Analysed Rotameric		Percentiles	
1	A	166/135 (123%)	151 (91%)	15 (9%)	9 2	

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	130[A]	ARG
1	A	148[A]	TRP
1	A	130[B]	ARG
1	A	148[B]	TRP
1	A	142[A]	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

6 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	Chain	Chain	Chain	Chain	Res	Link	Вс	Bond lengths			Bond angles		
MIOI	$oxed{egin{array}{c c} Mol & Type \\ \hline \end{array}}$	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2					
5	DIO	A	204	-	6,6,6	0.50	0	6,6,6	1.76	1 (16%)					
2	HEM	A	201	1	41,50,50	1.88	8 (19%)	45,82,82	2.32	17 (37%)					
4	SO4	A	203	-	4,4,4	0.16	0	6,6,6	0.30	0					
3	CMO	A	202	-	0,1,1	-	-	-							
5	DIO	A	205	-	6,6,6	0.50	0	6,6,6	0.88	0					
5	DIO	A	206	-	6,6,6	0.53	0	6,6,6	1.16	1 (16%)					

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
5	DIO	A	204	-	-	-	0/1/1/1
5	DIO	A	205	-	-	-	0/1/1/1
2	HEM	A	201	1	-	2/12/54/54	-
5	DIO	A	206	-	-	-	0/1/1/1

The worst 5 of 8 bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
2	A	201	HEM	C3D-C2D	7.01	1.51	1.36
2	A	201	HEM	C3C-CAC	4.08	1.56	1.47
2	A	201	HEM	FE-NB	3.06	2.12	1.96
2	A	201	HEM	C3C-C2C	-2.79	1.36	1.40
2	A	201	HEM	CAB-C3B	2.36	1.53	1.47

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	A	201	HEM	C4D-ND-C1D	7.48	112.80	105.07
2	A	201	HEM	C1B-NB-C4B	5.48	110.73	105.07
2	A	201	HEM	CHB-C1B-NB	4.58	130.04	124.38
2	A	201	HEM	C4B-C3B-C2B	4.44	110.64	107.11
2	A	201	HEM	CHD-C1D-ND	3.12	127.82	124.43

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	201	HEM	CAA-CBA-CGA-O2A
2	A	201	HEM	CAA-CBA-CGA-O1A

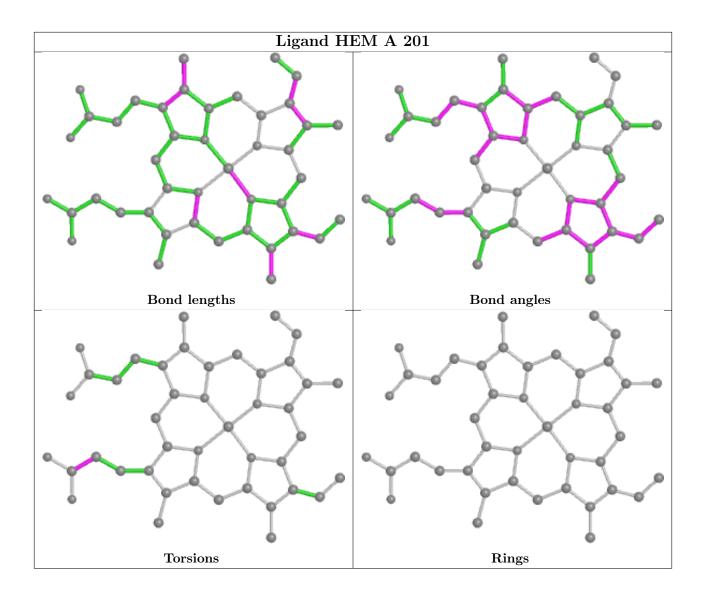
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	204	DIO	2	0
2	A	201	HEM	2	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 then 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 (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



6 Fit of model and data (i)

6.1 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, 95^{th} 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 $>$	$\# \mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q < 0.9
1	A	149/154 (96%)	0.27	13 (8%) 10 8	10, 21, 45, 72	0

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	ARG	6.4
1	A	148[A]	TRP	5.7
1	A	99[A]	VAL	5.4
1	A	98[A]	ALA	5.4
1	A	101[A]	VAL	5.2

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
5	DIO	A	204	6/6	0.82	0.24	27,29,34,35	0
5	DIO	A	206	6/6	0.90	0.34	13,19,25,26	6

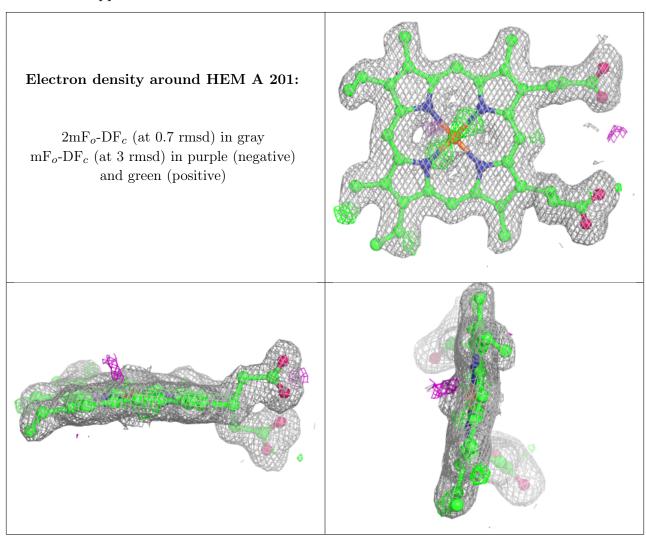
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
5	DIO	A	205	6/6	0.92	0.30	22,23,32,34	6
3	CMO	A	202	2/2	0.94	0.23	19,19,19,20	0
4	SO4	A	203	5/5	0.96	0.14	16,17,18,22	5
2	HEM	A	201	43/43	0.97	0.09	13,17,31,39	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.



6.5 Other polymers (i)

There are no such residues in this entry.

