



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 17, 2017 – 03:50 PM EDT

PDB ID : 5JGG
Title : X-ray sequence and high resolution crystal structure of Persian sturgeon methe-moglobin
Authors : Seyedarabi, A.
Deposited on : 2016-04-20
Resolution : 1.70 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029077
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029077

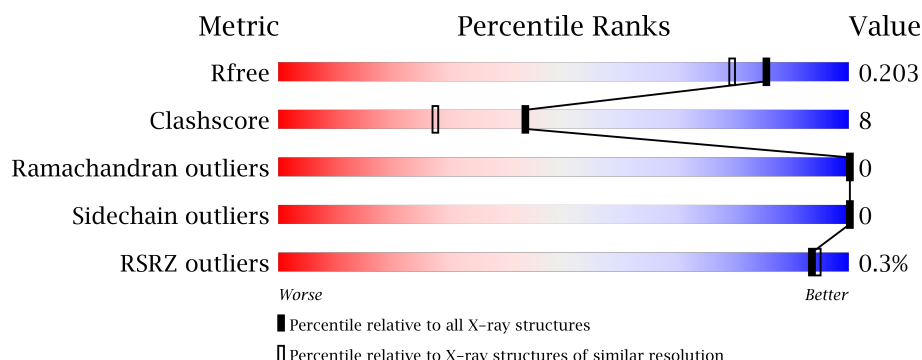
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	100719	3453 (1.70-1.70)
Clashscore	112137	3876 (1.70-1.70)
Ramachandran outliers	110173	3815 (1.70-1.70)
Sidechain outliers	110143	3815 (1.70-1.70)
RSRZ outliers	101464	3491 (1.70-1.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 $\leq 5\%$. 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	A	142	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0; width: 100%; text-align: center;">%</div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 15%, green 83%, grey 100%);"></div> <div style="position: absolute; bottom: -10px; left: 0; width: 100%; text-align: center;">83% 15% .</div> </div> </div>
2	B	146	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background: linear-gradient(to right, green 89%, yellow 10%, grey 100%);"></div> <div style="position: absolute; bottom: -10px; left: 0; width: 100%; text-align: center;">89% 10% .</div> </div> </div>

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
4	PGO	B	202	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2671 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.

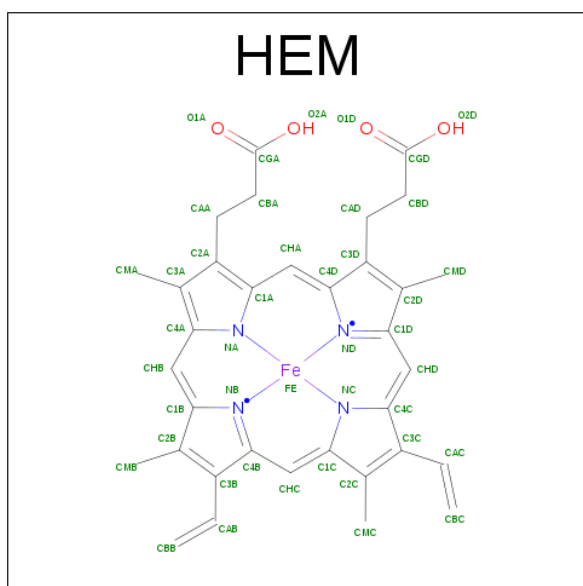
- Molecule 1 is a protein called Alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	142	Total	C	N	O	S	0	3	0
			1094	703	185	204	2			

- Molecule 2 is a protein called Beta chain.

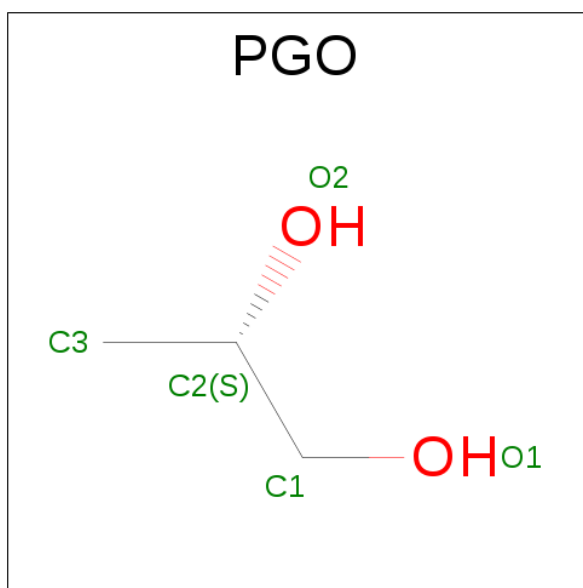
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	144	Total	C	N	O	S	0	1	0
			1106	710	184	209	3			

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
3	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 4 is S-1,2-PROPANEDIOL (three-letter code: PGO) (formula: $C_3H_8O_2$).

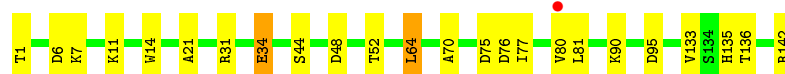


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			5	3	2		
4	B	1	Total	C	O	0	0
			5	3	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	199	Total	O	0	0
			199	199		
5	B	176	Total	O	0	0
			176	176		

- Molecule 1: Alpha chain



- Chain B: 89% 10%



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	53.89 Å 62.29 Å 101.56 Å 90.00° 97.67° 90.00°	Depositor
Resolution (Å)	52.97 – 1.70 33.26 – 1.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (52.97-1.70) 100.0 (33.26-1.70)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 1.70 Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.158 , 0.195 0.169 , 0.203	Depositor DCC
R_{free} test set	1802 reflections (5.16%)	DCC
Wilson B-factor (Å ²)	15.5	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2671	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.03% 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.

5 Model quality

5.1 Standard geometry

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

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	1.09	1/1139 (0.1%)	1.10	11/1550 (0.7%)
2	B	1.10	1/1139 (0.1%)	0.95	0/1559
All	All	1.09	2/2278 (0.1%)	1.03	11/3109 (0.4%)

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
1	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	34	GLU	CD-OE1	5.63	1.31	1.25
2	B	7	GLU	CD-OE2	-5.23	1.20	1.25

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	31	ARG	NE-CZ-NH1	8.29	124.45	120.30
1	A	142	ARG	CA-C-O	-7.74	103.84	120.10
1	A	48	ASP	CB-CG-OD1	6.56	124.21	118.30
1	A	142	ARG	NE-CZ-NH1	6.38	123.49	120.30
1	A	6	ASP	CB-CG-OD2	-6.11	112.80	118.30
1	A	48	ASP	CB-CG-OD2	-5.99	112.91	118.30
1	A	64	LEU	CA-CB-CG	-5.54	102.55	115.30
1	A	31	ARG	NE-CZ-NH2	-5.48	117.56	120.30
1	A	64	LEU	CB-CG-CD1	5.24	119.90	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	95	ASP	CB-CG-OD1	5.18	122.96	118.30
1	A	6	ASP	CB-CG-OD1	5.18	122.96	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	THR	Peptide

5.2 Too-close contacts ⓘ

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	1094	0	1107	21	1
2	B	1106	0	1064	10	1
3	A	43	0	30	0	0
3	B	43	0	30	5	0
4	B	10	0	16	0	0
5	A	199	0	0	8	0
5	B	176	0	0	4	0
All	All	2671	0	2247	35	1

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 8.

All (35) 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)
1:A:11:LYS:HE3	5:A:456:HOH:O	1.78	0.82
1:A:135:HIS:NE2	5:A:303:HOH:O	2.25	0.69
2:B:22:THR:HG23	5:B:304:HOH:O	1.92	0.69
2:B:22:THR:HG21	5:B:306:HOH:O	1.95	0.66
2:B:90:THR:HG21	5:B:327:HOH:O	1.95	0.66
1:A:21:ALA:HA	1:A:64:LEU:HD22	1.84	0.59
1:A:34:GLU:OE2	5:A:301:HOH:O	2.18	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:ALA:CB	1:A:64:LEU:HD22	2.36	0.56
1:A:77:ILE:HA	1:A:80:VAL:HG12	1.89	0.55
1:A:81:LEU:HB2	1:A:136:THR:CG2	2.37	0.55
1:A:80:VAL:HG13	1:A:81:LEU:HG	1.89	0.54
1:A:81:LEU:HB2	1:A:136:THR:HG21	1.88	0.54
1:A:81:LEU:CD1	1:A:133:VAL:HG13	2.38	0.53
1:A:7:LYS:NZ	1:A:75:ASP:OD1	2.30	0.52
3:B:201:HEM:HMC1	3:B:201:HEM:HBC2	1.92	0.52
1:A:135:HIS:CE1	5:A:303:HOH:O	2.62	0.52
2:B:94:GLU:OE1	5:B:301:HOH:O	2.19	0.51
2:B:21:GLU:HG2	2:B:65:LYS:HD2	1.92	0.51
1:A:76:ASP:OD1	5:A:302:HOH:O	2.19	0.50
1:A:90:LYS:NZ	5:A:308:HOH:O	2.42	0.50
3:B:201:HEM:CMC	3:B:201:HEM:HBC2	2.43	0.47
1:A:21:ALA:CA	1:A:64:LEU:HD22	2.45	0.47
1:A:44:SER:HA	5:A:433:HOH:O	2.14	0.46
2:B:25:ALA:O	2:B:29:VAL:HG23	2.15	0.46
1:A:52:THR:HG22	5:A:468:HOH:O	2.14	0.46
3:B:201:HEM:HBB2	3:B:201:HEM:CMB	2.46	0.45
2:B:8:ARG:O	2:B:12:VAL:HG22	2.16	0.45
3:B:201:HEM:HMB1	3:B:201:HEM:HBB2	2.00	0.43
1:A:70:ALA:HA	1:A:80:VAL:HG11	2.01	0.42
2:B:9:PHE:O	2:B:13[B]:THR:OG1	2.21	0.41
2:B:96:LEU:HD13	3:B:201:HEM:C2D	2.56	0.40
2:B:9:PHE:CZ	2:B:13[B]:THR:HG21	2.56	0.40

All (1) 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	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:SER:OG	2:B:59:GLN:NE2[1_455]	2.13	0.07

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	143/142 (101%)	142 (99%)	1 (1%)	0	100	100
2	B	143/146 (98%)	139 (97%)	4 (3%)	0	100	100
All	All	286/288 (99%)	281 (98%)	5 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	123/120 (102%)	123 (100%)	0	100	100
2	B	114/113 (101%)	114 (100%)	0	100	100
All	All	237/233 (102%)	237 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

4 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$
3	HEM	A	201	1,5	28,50,50	1.61	5 (17%)	17,82,82	2.15	7 (41%)
3	HEM	B	201	2,5	28,50,50	1.25	3 (10%)	17,82,82	1.85	5 (29%)
4	PGO	B	202	-	4,4,4	0.88	0	2,4,4	1.39	0
4	PGO	B	203	-	4,4,4	0.84	0	2,4,4	2.46	1 (50%)

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
3	HEM	A	201	1,5	-	0/6/54/54	0/0/8/8
3	HEM	B	201	2,5	-	0/6/54/54	0/0/8/8
4	PGO	B	202	-	-	0/2/2/2	0/0/0/0
4	PGO	B	203	-	-	0/2/2/2	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	201	HEM	C1B-NB	-3.61	1.32	1.36
3	A	201	HEM	C1C-NC	-3.08	1.33	1.36
3	B	201	HEM	C3B-C2B	-2.03	1.37	1.40
3	A	201	HEM	CMA-C3A	2.27	1.56	1.51
3	A	201	HEM	C4D-ND	2.48	1.39	1.36
3	B	201	HEM	C1C-NC	2.53	1.39	1.36
3	B	201	HEM	C4D-ND	3.52	1.40	1.36
3	A	201	HEM	C4C-NC	4.43	1.42	1.36

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	201	HEM	CBD-CAD-C3D	-4.40	104.08	112.47
3	A	201	HEM	CBD-CAD-C3D	-3.93	104.97	112.47
3	A	201	HEM	CMD-C2D-C1D	-3.22	123.51	128.46
3	A	201	HEM	CAA-CBA-CGA	-2.46	108.45	112.66
3	B	201	HEM	CMA-C3A-C4A	-2.35	124.86	128.46
3	A	201	HEM	CMA-C3A-C4A	-2.14	125.18	128.46
3	B	201	HEM	C3C-C4C-NC	-2.04	107.10	110.94
3	A	201	HEM	CMA-C3A-C2A	2.12	128.95	124.94
3	A	201	HEM	CMD-C2D-C3D	2.32	129.31	124.94
3	B	201	HEM	CAD-CBD-CGD	2.55	117.02	112.66
3	B	201	HEM	CMC-C2C-C3C	3.02	130.49	124.89
4	B	203	PGO	O1-C1-C2	3.48	118.71	110.83
3	A	201	HEM	CAD-CBD-CGD	3.50	118.64	112.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	201	HEM	5	0

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, 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	142/142 (100%)	-0.34	1 (0%) 87 90	8, 16, 32, 39	0
2	B	144/146 (98%)	-0.30	0 100 100	8, 16, 33, 42	0
All	All	286/288 (99%)	-0.32	1 (0%) 93 94	8, 16, 33, 42	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	80	VAL	2.1

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	PGO	B	202	5/5	0.93	0.14	2.23	31,33,35,39	0
4	PGO	B	203	5/5	0.81	0.14	1.72	31,36,40,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	HEM	A	201	43/43	0.98	0.07	0.08	9,12,26,45	0
3	HEM	B	201	43/43	0.98	0.07	0.03	12,14,31,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.