



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 11:12 pm GMT

PDB ID : 1FSL  
Title : FERRIC SOYBEAN LEGHEMOGLOBIN COMPLEXED WITH NICOTINATE  
Authors : Ellis, P.J.; Guss, J.M.; Freeman, H.C.  
Deposited on : 1995-12-12  
Resolution : 2.30 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](mailto: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.

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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	:	trunk28620
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)	:	recalc28949

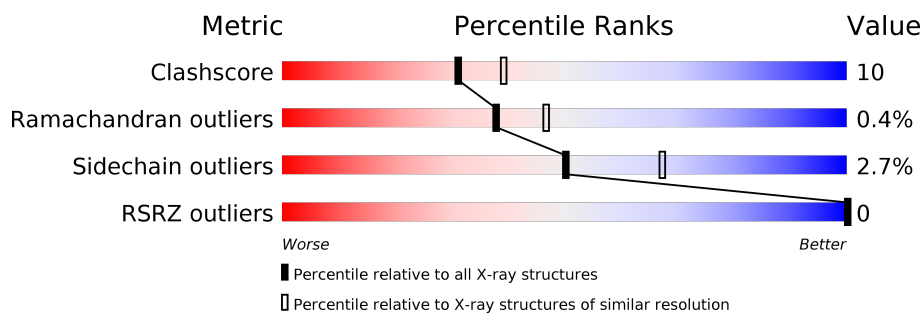
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*



The reported resolution of this entry is 2.30 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))
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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  $\geq 3$ , 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	143	 79% 19% ..
1	B	143	 66% 31% ..

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4635 atoms, of which 2293 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 LEGHEMOGLOBIN A.

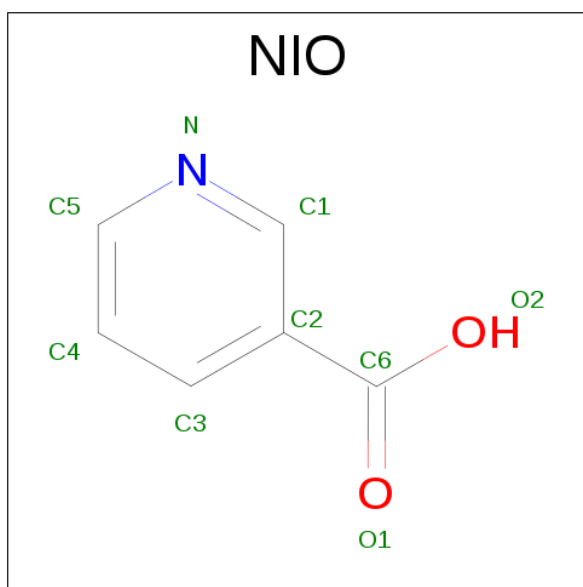
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	143	Total	C	H	N	O	0	2	0
			2195	698	1111	179	207			
1	B	143	Total	C	H	N	O	0	3	0
			2198	697	1114	180	207			

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	H	N	O	0
			73	34	1	30	4	4	
2	B	1	Total	C	Fe	H	N	O	0
			73	34	1	30	4	4	

- Molecule 3 is NICOTINIC ACID (three-letter code: NIO) (formula:  $C_6H_5NO_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			13	6	4	1	2		
3	B	1	Total	C	H	N	O	0	0
			13	6	4	1	2		


- Molecule 4 is water.

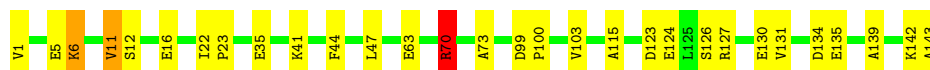
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	39	Total	O	0	0
			39	39		
4	B	31	Total	O	0	0
			31	31		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: LEGHEMOGLOBIN A

Chain A:  79% 19% ..



#### • Molecule 1: LEGHEMOGLOBIN A

Chain B:  66% 31% ..



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.69Å 56.55Å 57.80Å 90.00° 109.90° 90.00°	Depositor
Resolution (Å)	6.00 – 2.30 39.19 – 2.31	Depositor EDS
% Data completeness (in resolution range)	49.5 (6.00-2.30) 53.3 (39.19-2.31)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	31.24 (at 2.31Å)	Xtriage
Refinement program	PROFFT	Depositor
R, $R_{free}$	0.158 , (Not available) 0.161 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	12.9	Xtriage
Anisotropy	0.323	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 56.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.38$ , $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4635	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.53% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.92	4/1118 (0.4%)	1.10	4/1514 (0.3%)
1	B	0.94	7/1127 (0.6%)	1.10	5/1525 (0.3%)
All	All	0.93	11/2245 (0.5%)	1.10	9/3039 (0.3%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	63	GLU	CD-OE2	6.62	1.32	1.25
1	B	63	GLU	CD-OE2	6.60	1.32	1.25
1	A	5	GLU	CD-OE1	6.19	1.32	1.25
1	B	35	GLU	CD-OE1	6.16	1.32	1.25
1	B	124	GLU	CD-OE1	6.01	1.32	1.25
1	A	16	GLU	CD-OE1	5.91	1.32	1.25
1	B	5	GLU	CD-OE1	5.90	1.32	1.25
1	A	35	GLU	CD-OE1	5.63	1.31	1.25
1	B	130	GLU	CD-OE1	5.49	1.31	1.25
1	B	107	GLU	CD-OE2	5.47	1.31	1.25
1	B	16	GLU	CD-OE1	5.20	1.31	1.25

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	134	ASP	CB-CG-OD1	8.11	125.60	118.30
1	A	123	ASP	CB-CG-OD1	6.13	123.81	118.30
1	B	70	ARG	NE-CZ-NH1	6.09	123.35	120.30
1	A	134	ASP	CB-CG-OD1	5.95	123.66	118.30
1	A	70	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	B	134	ASP	CB-CG-OD2	-5.85	113.04	118.30
1	B	127	ARG	NE-CZ-NH1	5.59	123.09	120.30
1	B	85	ASP	CB-CG-OD1	5.26	123.04	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	123	ASP	CB-CG-OD2	-5.20	113.62	118.30

There are no chirality outliers.

There are no planarity outliers.

## 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	1084	1111	1100	17	0
1	B	1084	1114	1093	29	0
2	A	43	30	30	1	0
2	B	43	30	30	1	0
3	A	9	4	4	1	0
3	B	9	4	4	0	0
4	A	39	0	0	2	0
4	B	31	0	0	0	0
All	All	2342	2293	2261	47	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 10.

All (47) 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:139:ALA:O	1:A:143:ALA:HB3	1.89	0.71
1:A:126:SER:O	1:A:130:GLU:HB2	1.95	0.66
1:B:38:PRO:O	1:B:41:LYS:HE2	1.94	0.66
1:B:34:LEU:HD23	1:B:40:ALA:HB3	1.78	0.65
1:B:127:ARG:O	1:B:131:VAL:HG23	1.96	0.64
1:A:44:PHE:HB2	1:A:47:LEU:HD12	1.80	0.62
1:B:123:ASP:O	1:B:127:ARG:HG3	2.01	0.61
1:A:1:VAL:HG21	1:A:135:GLU:OE1	2.01	0.60
1:B:4:THR:OG1	1:B:7:GLN:HG3	2.01	0.60
1:A:22:ILE:N	1:A:23:PRO:CD	2.66	0.59
1:A:127:ARG:O	1:A:131:VAL:HG23	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:ALA:O	1:B:143:ALA:HB3	2.05	0.57
1:B:9:ALA:O	1:B:13:SER:HB3	2.06	0.56
1:B:6:LYS:HE3	1:B:6:LYS:HA	1.89	0.55
1:B:100:PRO:O	1:B:104:VAL:HG23	2.07	0.54
1:B:12:SER:HB3	1:B:70:ARG:NH2	2.23	0.54
1:B:6:LYS:HA	1:B:6:LYS:CE	2.37	0.53
1:B:99:ASP:N	1:B:100:PRO:CD	2.73	0.51
1:B:141:LYS:O	1:B:142:LYS:CB	2.58	0.51
1:B:142:LYS:O	1:B:143:ALA:HB2	2.10	0.50
1:B:142:LYS:O	1:B:143:ALA:CB	2.60	0.50
1:B:22:ILE:N	1:B:23:PRO:CD	2.75	0.50
1:B:3:PHE:CE2	1:B:73:ALA:HB1	2.48	0.48
1:A:41:LYS:HE2	1:A:47:LEU:O	2.13	0.48
1:A:99:ASP:N	1:A:100:PRO:CD	2.76	0.48
1:B:141:LYS:O	1:B:142:LYS:HB3	2.13	0.48
1:A:11:VAL:HG11	1:A:73:ALA:CB	2.45	0.46
1:B:92:HIS:HA	1:B:97:VAL:HG21	1.97	0.46
1:B:84:ALA:HB1	1:B:88:LEU:HD23	1.98	0.45
1:A:142:LYS:O	1:A:143:ALA:HB2	2.17	0.45
1:A:12:SER:OG	1:A:70:ARG:NH2	2.51	0.44
1:B:22:ILE:N	1:B:23:PRO:HD2	2.32	0.44
1:A:131:VAL:HG13	4:A:164:HOH:O	2.17	0.44
1:B:68:LEU:O	1:B:71:ASP:HB3	2.18	0.43
1:B:29:PHE:CZ	1:B:33:ILE:HD11	2.54	0.43
1:B:88:LEU:O	1:B:92:HIS:ND1	2.42	0.43
1:A:6[B]:LYS:HE3	1:A:124:GLU:OE2	2.19	0.42
1:B:22:ILE:HB	1:B:23:PRO:HD3	2.00	0.42
1:B:51:VAL:O	1:B:53:PRO:HD3	2.20	0.42
1:B:121:TRP:CE3	1:B:125:LEU:HD23	2.55	0.42
1:B:98:THR:OG1	1:B:100:PRO:HD2	2.20	0.42
2:A:144:HEM:NA	3:A:145:NIO:H1	2.35	0.41
1:A:115:ALA:HB2	4:A:163:HOH:O	2.21	0.41
1:B:140:ILE:HD12	2:B:144:HEM:HBB2	2.02	0.41
1:A:99:ASP:HB2	1:A:100:PRO:HD3	2.03	0.40

There are no symmetry-related clashes.

## 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	143/143 (100%)	138 (96%)	5 (4%)	0	100	100
1	B	144/143 (101%)	138 (96%)	5 (4%)	1 (1%)	25	30
All	All	287/286 (100%)	276 (96%)	10 (4%)	1 (0%)	38	55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	142	LYS

### 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	113/111 (102%)	109 (96%)	4 (4%)	41	56
1	B	114/111 (103%)	111 (97%)	3 (3%)	51	69
All	All	227/222 (102%)	220 (97%)	7 (3%)	50	61

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6[A]	LYS
1	A	6[B]	LYS
1	A	11	VAL
1	A	70	ARG
1	B	6	LYS

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Mol	Chain	Res	Type
1	B	70	ARG
1	B	119	ASP

Some 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	$\# Z  > 2$
2	HEM	A	144	1,3	28,50,50	1.96	7 (25%)	17,82,82	2.16	9 (52%)
3	NIO	A	145	2	6,9,9	1.16	1 (16%)	8,11,11	1.22	1 (12%)
2	HEM	B	144	1,3	28,50,50	2.01	7 (25%)	17,82,82	1.86	5 (29%)
3	NIO	B	145	2	6,9,9	0.95	0	8,11,11	1.34	1 (12%)

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	HEM	A	144	1,3	-	0/6/54/54	0/0/8/8
3	NIO	A	145	2	-	0/0/4/4	0/1/1/1
2	HEM	B	144	1,3	-	0/6/54/54	0/0/8/8
3	NIO	B	145	2	-	0/0/4/4	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	144	HEM	C3B-C2B	-5.45	1.33	1.40
2	A	144	HEM	C3B-C2B	-5.29	1.33	1.40
2	A	144	HEM	C3C-C2C	-4.27	1.34	1.40
2	B	144	HEM	C3C-C2C	-4.04	1.35	1.40
2	A	144	HEM	C1B-NB	2.02	1.39	1.36
2	A	144	HEM	CAA-C2A	2.06	1.55	1.52
2	B	144	HEM	CAA-C2A	2.07	1.55	1.52
2	A	144	HEM	C4C-NC	2.23	1.39	1.36
2	B	144	HEM	C4C-NC	2.25	1.39	1.36
3	A	145	NIO	C1-N	2.41	1.39	1.34
2	B	144	HEM	C1C-NC	2.49	1.39	1.36
2	B	144	HEM	C3C-CAC	3.12	1.53	1.47
2	A	144	HEM	C3B-CAB	3.13	1.54	1.47
2	A	144	HEM	C3C-CAC	3.46	1.54	1.47
2	B	144	HEM	C3B-CAB	3.64	1.55	1.47

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	144	HEM	CMD-C2D-C1D	-4.15	122.09	128.46
2	A	144	HEM	CMA-C3A-C4A	-3.43	123.19	128.46
2	B	144	HEM	CMD-C2D-C1D	-2.93	123.97	128.46
2	B	144	HEM	CMA-C3A-C4A	-2.89	124.02	128.46
2	A	144	HEM	CMB-C2B-C3B	2.05	128.69	124.89
2	A	144	HEM	CMC-C2C-C3C	2.16	128.90	124.89
2	A	144	HEM	CMA-C3A-C2A	2.21	129.11	124.94
3	B	145	NIO	C3-C4-C5	2.24	122.24	118.91
3	A	145	NIO	C3-C4-C5	2.25	122.26	118.91
2	A	144	HEM	CBD-CAD-C3D	2.32	116.90	112.47
2	B	144	HEM	CMB-C2B-C3B	2.38	129.31	124.89
2	B	144	HEM	CAA-CBA-CGA	2.64	117.17	112.66
2	A	144	HEM	CBA-CAA-C2A	2.86	117.95	112.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	144	HEM	CMD-C2D-C3D	2.99	130.59	124.94
2	A	144	HEM	CAA-CBA-CGA	3.00	117.79	112.66
2	B	144	HEM	CBA-CAA-C2A	3.05	118.31	112.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	144	HEM	1	0
3	A	145	NIO	1	0
2	B	144	HEM	1	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	143/143 (100%)	-0.51	0 100 100	2, 12, 21, 31	0
1	B	143/143 (100%)	-0.50	0 100 100	3, 12, 22, 28	0
All	All	286/286 (100%)	-0.51	0 100 100	2, 12, 22, 31	0

There are no RSRZ outliers to report.

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	HEM	A	144	43/43	0.97	0.13	1.20	12,14,16,17	0
2	HEM	B	144	43/43	0.97	0.12	0.90	8,10,15,16	0
3	NIO	A	145	9/9	0.98	0.11	-0.06	5,6,6,6	0
3	NIO	B	145	9/9	0.97	0.10	-0.24	4,5,6,6	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.