



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 02:44 am GMT

PDB ID : 1DGB
Title : HUMAN ERYTHROCYTE CATALASE
Authors : Putnam, C.D.; Arvai, A.S.; Bourne, Y.; Tainer, J.A.
Deposited on : 1999-11-23
Resolution : 2.20 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

<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 : 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

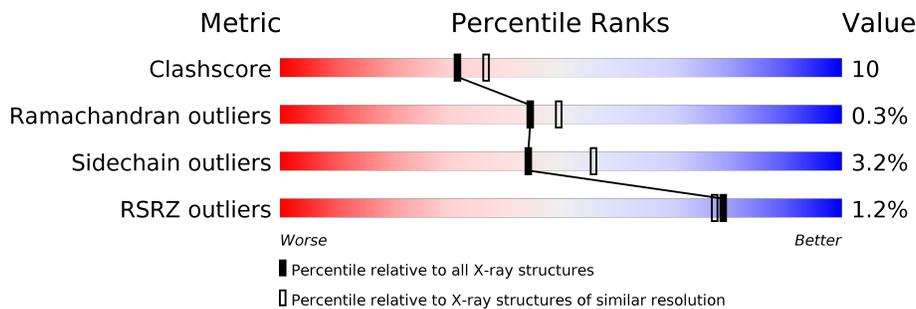
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

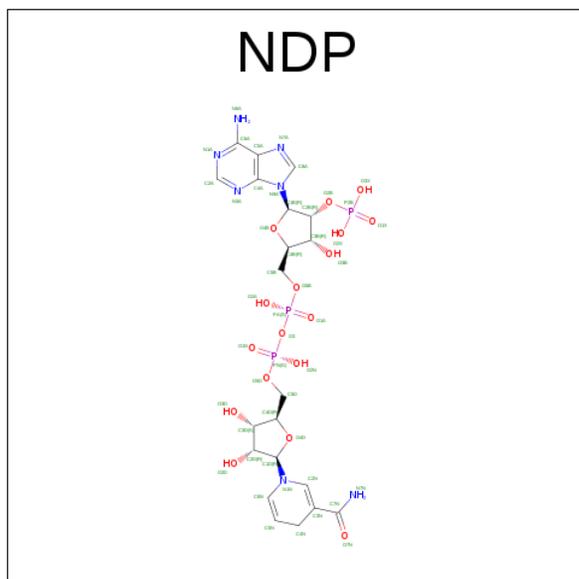
The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey 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	498	 % 83% 16% .
1	B	498	 % 77% 22% .
1	C	498	 % 77% 21% .
1	D	498	 % 77% 21% .

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

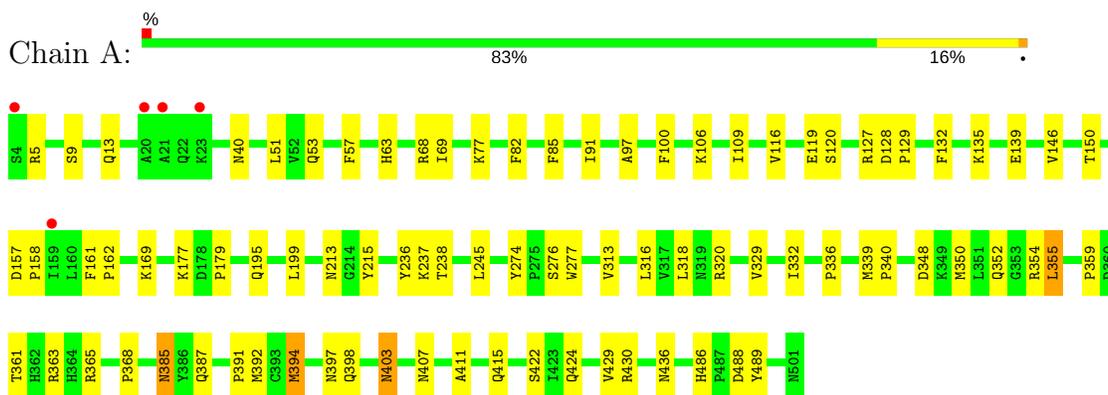
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	309	Total	O	0	0
			309	309		
4	B	263	Total	O	0	0
			263	263		
4	C	266	Total	O	0	0
			266	266		
4	D	271	Total	O	0	0
			271	271		

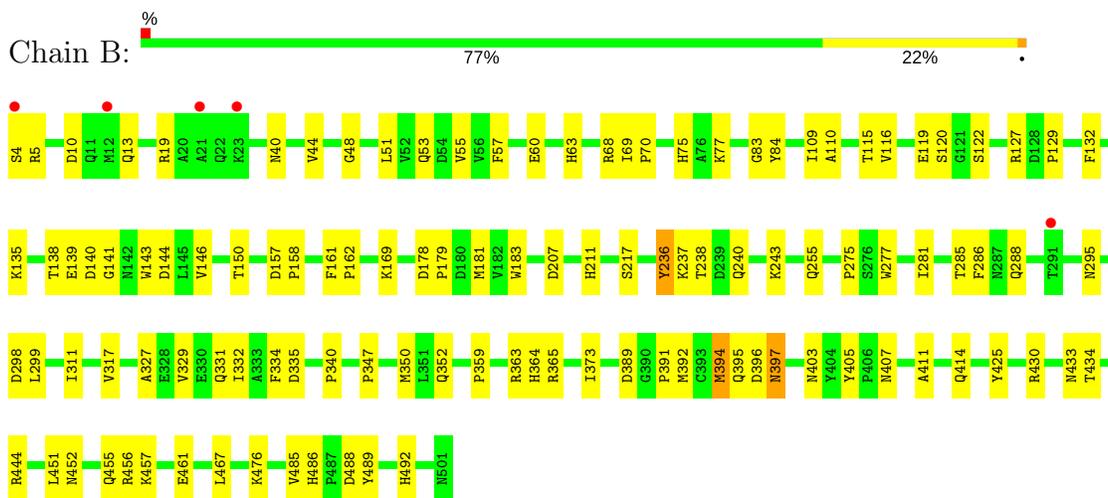
3 Residue-property plots [i](#)

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.

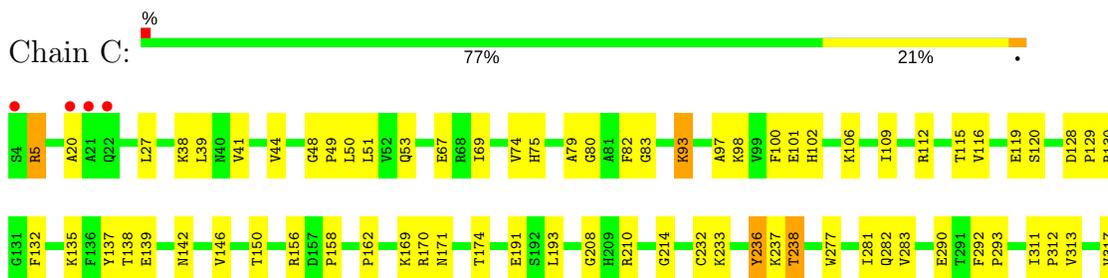
- Molecule 1: CATALASE



- Molecule 1: CATALASE

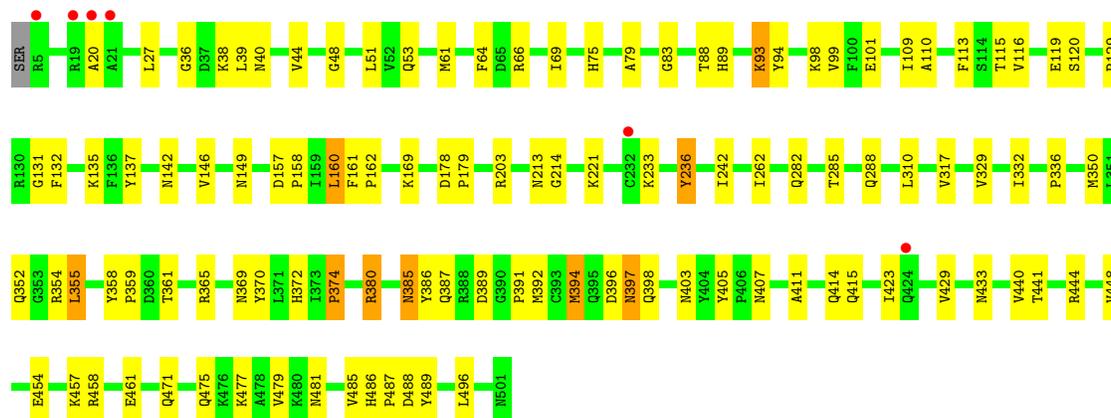
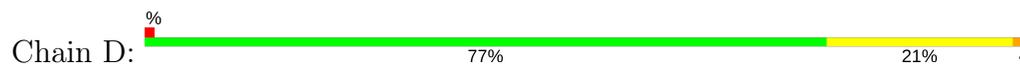


- Molecule 1: CATALASE





● Molecule 1: CATALASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.50Å 139.93Å 227.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 19.96 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.20) 90.6 (19.96-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.11 (at 2.19Å)	Xtrriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.172 , 0.227 0.172 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	64.4	Xtrriage
Anisotropy	0.051	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 55.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	17443	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.82% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.38	0/4137	0.62	0/5619
1	B	0.39	0/4137	0.62	0/5619
1	C	0.39	0/4137	0.64	0/5619
1	D	0.40	0/4131	0.63	0/5611
All	All	0.39	0/16542	0.63	0/22468

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	4018	0	3845	81	0
1	B	4018	0	3845	102	0
1	C	4018	0	3845	96	0
1	D	4012	0	3840	104	0
2	A	43	0	30	1	0
2	B	43	0	30	0	0
2	C	43	0	30	0	0
2	D	43	0	30	0	0
3	A	48	0	26	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	48	0	26	0	0
4	A	309	0	0	18	0
4	B	263	0	0	10	0
4	C	266	0	0	12	0
4	D	271	0	0	15	0
All	All	17443	0	15547	325	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.

The worst 5 of 325 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:B:138:THR:HG22	1:B:140:ASP:H	1.17	1.05
1:B:452:ASN:H	1:B:455:GLN:HE21	0.99	0.92
1:A:13:GLN:HE21	1:C:471:GLN:HE22	1.12	0.92
1:A:120[B]:SER:O	1:C:120[B]:SER:O	1.89	0.91
1:A:53:GLN:HE21	1:D:352:GLN:HE22	1.19	0.89

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	498/498 (100%)	463 (93%)	35 (7%)	0	100	100
1	B	498/498 (100%)	468 (94%)	30 (6%)	0	100	100
1	C	498/498 (100%)	466 (94%)	30 (6%)	2 (0%)	38	41
1	D	497/498 (100%)	467 (94%)	27 (5%)	3 (1%)	28	29
All	All	1991/1992 (100%)	1864 (94%)	122 (6%)	5 (0%)	44	49

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	396	ASP
1	C	20	ALA
1	D	20	ALA
1	C	441	THR
1	D	374	PRO

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	430/429 (100%)	416 (97%)	14 (3%)	43 54
1	B	430/429 (100%)	419 (97%)	11 (3%)	51 64
1	C	430/429 (100%)	413 (96%)	17 (4%)	36 45
1	D	429/429 (100%)	416 (97%)	13 (3%)	46 58
All	All	1719/1716 (100%)	1664 (97%)	55 (3%)	44 56

5 of 55 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	492	HIS
1	C	138	THR
1	D	385	ASN
1	C	5	ARG
1	C	67	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 59 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	475	GLN
1	C	338	ASN
1	D	403	ASN
1	B	486	HIS
1	C	40	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 carbohydrates 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	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HEM	A	1200	1	28,50,50	1.78	6 (21%)	17,82,82	2.03	6 (35%)
3	NDP	A	1300	-	43,52,52	1.14	2 (4%)	49,80,80	1.51	5 (10%)
2	HEM	B	1201	1	28,50,50	1.82	4 (14%)	17,82,82	1.89	6 (35%)
2	HEM	C	1202	1	28,50,50	1.71	5 (17%)	17,82,82	2.05	5 (29%)
3	NDP	C	1301	-	43,52,52	1.15	2 (4%)	49,80,80	1.40	3 (6%)
2	HEM	D	1203	1	28,50,50	1.81	5 (17%)	17,82,82	2.01	5 (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	HEM	A	1200	1	-	0/6/54/54	0/0/8/8
3	NDP	A	1300	-	-	0/30/77/77	0/5/5/5
2	HEM	B	1201	1	-	0/6/54/54	0/0/8/8

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	C	1202	1	-	0/6/54/54	0/0/8/8
3	NDP	C	1301	-	-	0/30/77/77	0/5/5/5
2	HEM	D	1203	1	-	0/6/54/54	0/0/8/8

The worst 5 of 24 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1201	HEM	C3C-C2C	-6.19	1.32	1.40
2	C	1202	HEM	C3B-C2B	-5.48	1.33	1.40
2	A	1200	HEM	C3B-C2B	-5.31	1.33	1.40
2	D	1203	HEM	C3B-C2B	-5.08	1.33	1.40
2	D	1203	HEM	C3C-C2C	-4.85	1.34	1.40

The worst 5 of 30 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1300	NDP	N3A-C2A-N1A	-6.76	122.97	128.86
3	C	1301	NDP	N3A-C2A-N1A	-6.59	123.12	128.86
2	C	1202	HEM	CMD-C2D-C1D	-5.08	120.66	128.46
2	A	1200	HEM	CMA-C3A-C4A	-4.20	122.01	128.46
2	D	1203	HEM	CMD-C2D-C1D	-3.26	123.45	128.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1200	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, 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	498/498 (100%)	-0.33	5 (1%) 82 81	55, 62, 71, 83	0
1	B	498/498 (100%)	-0.34	5 (1%) 82 81	55, 62, 72, 82	0
1	C	498/498 (100%)	-0.35	7 (1%) 75 73	54, 62, 74, 86	0
1	D	497/498 (99%)	-0.31	6 (1%) 79 77	54, 61, 73, 85	0
All	All	1991/1992 (99%)	-0.33	23 (1%) 79 77	54, 62, 73, 86	0

The worst 5 of 23 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	4	SER	5.9
1	B	4	SER	5.7
1	A	4	SER	5.2
1	D	21	ALA	5.2
1	A	21	ALA	4.6

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(\AA^2)	Q < 0.9
2	HEM	D	1203	43/43	0.97	0.12	-0.13	54,56,56,57	0
3	NDP	C	1301	48/48	0.95	0.11	-0.43	73,79,83,83	0
2	HEM	B	1201	43/43	0.98	0.09	-0.61	56,57,58,59	0
2	HEM	C	1202	43/43	0.97	0.10	-0.62	58,60,61,61	0
2	HEM	A	1200	43/43	0.98	0.09	-0.73	56,57,58,58	0
3	NDP	A	1300	48/48	0.96	0.09	-0.74	67,71,74,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.