



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 08:46 PM GMT

PDB ID : 1LY8
Title : The crystal structure of a mutant enzyme of Coprinus cinereus peroxidase provides an understanding of its increased thermostability and insight into modelling of protein structures
Authors : Houborg, K.; Harris, P.; Poulsen, J.-C.N.; Svendsen, A.; Schneider, P.; Larsen, S.
Deposited on : 2002-06-07
Resolution : 2.05 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 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
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)	:	trunk26865

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BMA	B	600	-	-	-	X
7	GOL	A	801	-	-	-	X
7	GOL	A	806	-	-	-	X
7	GOL	A	810	-	-	-	X
7	GOL	A	811	-	-	-	X
7	GOL	B	803	-	-	-	X
7	GOL	B	804	-	-	-	X
7	GOL	B	805	-	-	-	X
7	GOL	B	807	-	-	-	X
7	GOL	B	808	-	-	-	X
7	GOL	B	809	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5874 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 Peroxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	339	Total	C	N	O	S	0	5	0
			2504	1559	430	501	14			
1	B	338	Total	C	N	O	S	0	5	0
			2499	1555	429	501	14			

There are 14 discrepancies between the modelled and reference sequences:

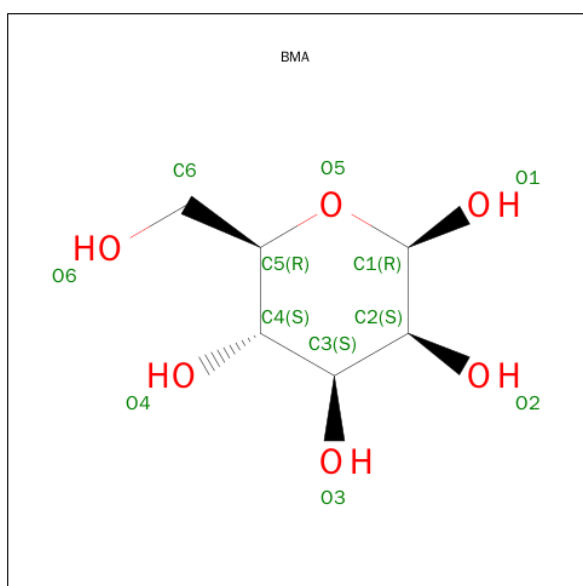
Chain	Residue	Modelled	Actual	Comment	Reference
A	49	SER	ILE	ENGINEERED	UNP P28314
A	53	ALA	VAL	ENGINEERED	UNP P28314
A	121	ALA	THR	ENGINEERED	UNP P28314
A	166	PHE	MET	ENGINEERED	UNP P28314
A	239	GLY	GLU	ENGINEERED	UNP P28314
A	242	ILE	MET	ENGINEERED	UNP P28314
A	272	PHE	TYR	ENGINEERED	UNP P28314
B	49	SER	ILE	ENGINEERED	UNP P28314
B	53	ALA	VAL	ENGINEERED	UNP P28314
B	121	ALA	THR	ENGINEERED	UNP P28314
B	166	PHE	MET	ENGINEERED	UNP P28314
B	239	GLY	GLU	ENGINEERED	UNP P28314
B	242	ILE	MET	ENGINEERED	UNP P28314
B	272	PHE	TYR	ENGINEERED	UNP P28314

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



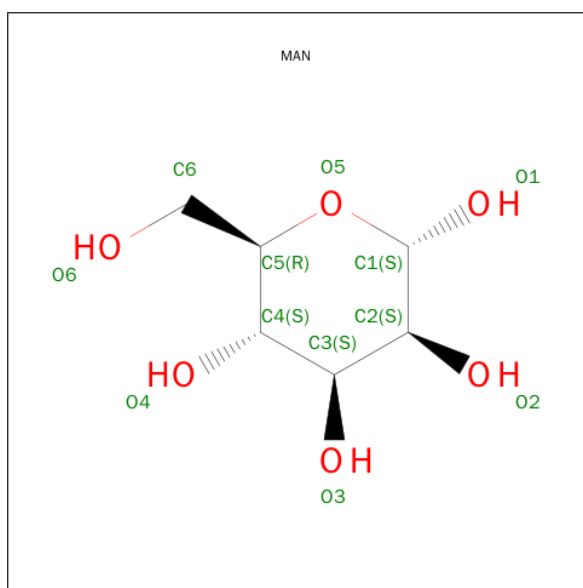
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is SUGAR (ALPHA-D-MANNOSE) (three-letter code: BMA) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			11	6	5		
3	B	1	Total	C	O	0	0
			11	6	5		

- Molecule 4 is SUGAR (ALPHA-D-MANNOSE) (three-letter code: MAN) (formula: $C_6H_{12}O_6$).

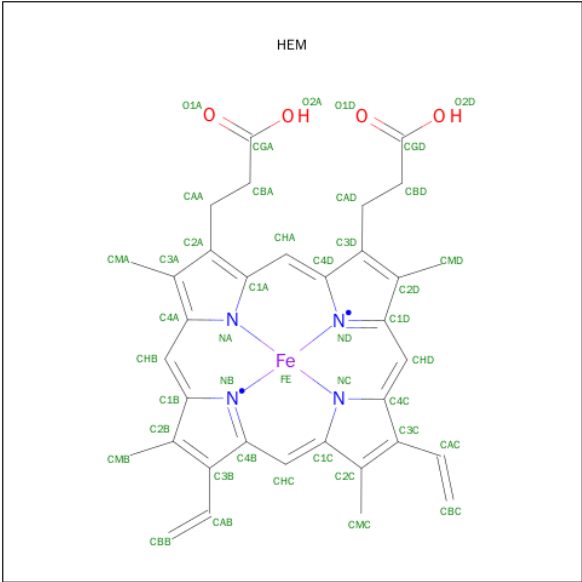


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			11	6	5		
4	B	1	Total	C	O	0	0
			11	6	5		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

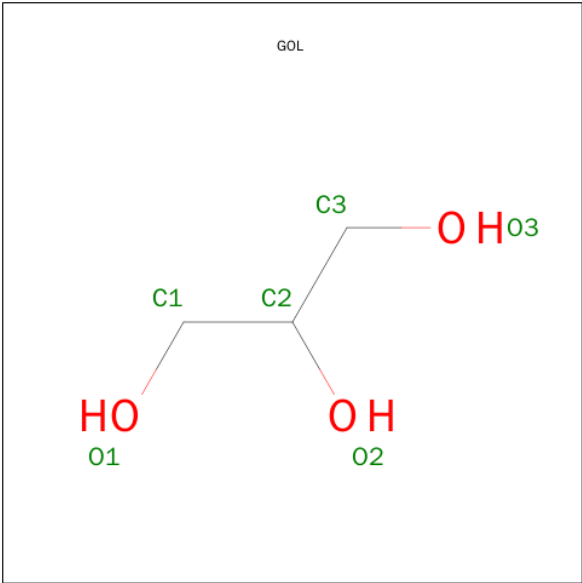
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Ca	0	0
			2	2		
5	A	2	Total	Ca	0	0
			2	2		

- Molecule 6 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		

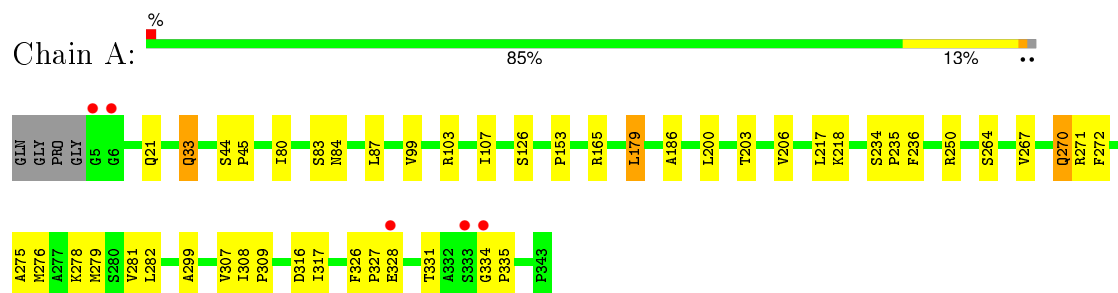
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	332	Total	O	0	0
			332	332		
8	B	305	Total	O	0	0
			305	305		

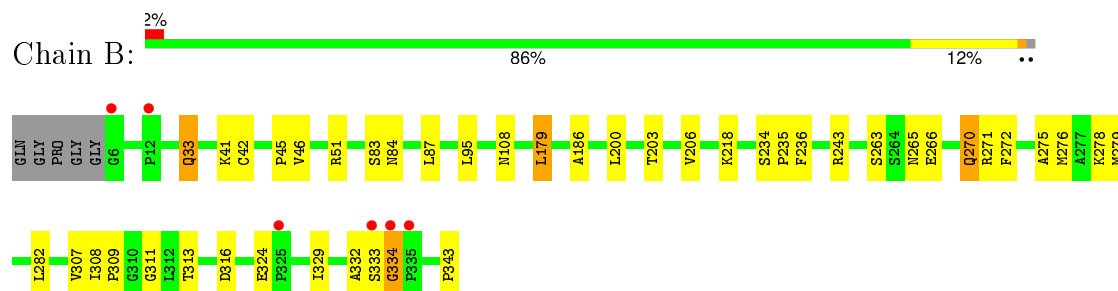
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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: Peroxidase



• Molecule 1: Peroxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	74.49Å 114.86Å 73.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.91 – 2.05 19.91 – 2.05	Depositor EDS
% Data completeness (in resolution range)	84.6 (19.91-2.05) 84.8 (19.91-2.05)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.39 (at 2.06Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.202 , 0.249 0.200 , 0.247	Depositor DCC
R_{free} test set	1687 reflections (5.22%)	DCC
Wilson B-factor (Å ²)	13.3	Xtriage
Anisotropy	0.869	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 61.0	EDS
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Outliers	1 of 33983 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5874	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.86 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.9183e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, BMA, NAG, CA, HEM, MAN

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	0.31	0/2561	0.58	0/3489
1	B	0.31	0/2556	0.57	0/3483
All	All	0.31	0/5117	0.57	0/6972

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	335	PRO	Mainchain

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	2504	0	2395	34	0
1	B	2499	0	2388	35	0
2	A	14	0	13	0	0
2	B	14	0	13	0	0
3	A	11	0	10	1	0
3	B	11	0	10	1	0
4	A	11	0	10	1	0
4	B	11	0	10	0	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	43	0	30	0	0
6	B	43	0	30	0	0
7	A	24	0	28	0	0
7	B	48	0	56	3	0
8	A	332	0	0	3	0
8	B	305	0	0	3	0
All	All	5874	0	4993	70	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 7.

The worst 5 of 70 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:33:GLN:HA	1:A:33:GLN:HE21	1.39	0.87
1:B:33:GLN:HE21	1:B:33:GLN:HA	1.44	0.83
1:B:265:ASN:H	7:B:808:GOL:H2	1.56	0.70
1:A:179:LEU:HD13	1:A:275:ALA:HB1	1.77	0.67
1:B:236:PHE:CZ	1:B:309:PRO:HG2	2.32	0.64

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	342/343 (100%)	328 (96%)	13 (4%)	1 (0%)	46	36
1	B	341/343 (99%)	332 (97%)	8 (2%)	1 (0%)	46	36
All	All	683/686 (100%)	660 (97%)	21 (3%)	2 (0%)	46	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	334	GLY
1	A	334	GLY

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	273/270 (101%)	268 (98%)	5 (2%)	66	62
1	B	273/270 (101%)	269 (98%)	4 (2%)	72	70
All	All	546/540 (101%)	537 (98%)	9 (2%)	68	67

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	328	GLU
1	B	270	GLN
1	B	84	ASN
1	A	179	LEU
1	B	33	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	302	ASN
1	B	21	GLN
1	B	127	ASN
1	A	270	GLN

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Mol	Chain	Res	Type
1	B	69	GLN

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 ⓘ

Of 24 ligands modelled in this entry, 4 are monoatomic - leaving 20 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	$\# Z > 2$
6	HEM	A	344	1	30,50,50	3.43	20 (66%)	24,82,82	4.44	13 (54%)
2	NAG	A	500	1	14,14,15	0.56	0	15,19,21	0.85	1 (6%)
3	BMA	A	600	1	11,11,12	0.50	0	14,15,17	0.63	0
4	MAN	A	700	1	11,11,12	0.41	0	14,15,17	0.43	0
7	GOL	A	801	-	5,5,5	1.04	0	5,5,5	1.76	2 (40%)
7	GOL	A	806	-	5,5,5	1.07	0	5,5,5	1.76	1 (20%)
7	GOL	A	810	-	5,5,5	1.15	0	5,5,5	1.76	1 (20%)
7	GOL	A	811	-	5,5,5	1.11	0	5,5,5	1.63	1 (20%)
6	HEM	B	344	1	30,50,50	3.97	18 (60%)	24,82,82	3.56	12 (50%)
2	NAG	B	500	1	14,14,15	0.59	0	15,19,21	0.84	1 (6%)
3	BMA	B	600	1	11,11,12	0.45	0	14,15,17	0.63	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MAN	B	700	1	11,11,12	0.43	0	14,15,17	0.44	0
7	GOL	B	800	-	5,5,5	1.08	0	5,5,5	1.70	2 (40%)
7	GOL	B	802	-	5,5,5	1.09	0	5,5,5	1.73	2 (40%)
7	GOL	B	803	-	5,5,5	1.11	0	5,5,5	1.78	2 (40%)
7	GOL	B	804	-	5,5,5	1.06	0	5,5,5	1.72	2 (40%)
7	GOL	B	805	-	5,5,5	1.04	0	5,5,5	1.70	1 (20%)
7	GOL	B	807	-	5,5,5	1.08	0	5,5,5	1.69	1 (20%)
7	GOL	B	808	-	5,5,5	1.05	0	5,5,5	1.74	2 (40%)
7	GOL	B	809	-	5,5,5	1.07	0	5,5,5	1.70	1 (20%)

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
6	HEM	A	344	1	-	0/10/54/54	0/0/8/8
2	NAG	A	500	1	-	0/6/23/26	0/1/1/1
3	BMA	A	600	1	-	0/2/19/22	0/1/1/1
4	MAN	A	700	1	-	0/2/19/22	0/1/1/1
7	GOL	A	801	-	-	0/4/4/4	0/0/0/0
7	GOL	A	806	-	-	0/4/4/4	0/0/0/0
7	GOL	A	810	-	-	0/4/4/4	0/0/0/0
7	GOL	A	811	-	-	0/4/4/4	0/0/0/0
6	HEM	B	344	1	-	0/10/54/54	0/0/8/8
2	NAG	B	500	1	-	0/6/23/26	0/1/1/1
3	BMA	B	600	1	-	0/2/19/22	0/1/1/1
4	MAN	B	700	1	-	0/2/19/22	0/1/1/1
7	GOL	B	800	-	-	0/4/4/4	0/0/0/0
7	GOL	B	802	-	-	0/4/4/4	0/0/0/0
7	GOL	B	803	-	-	0/4/4/4	0/0/0/0
7	GOL	B	804	-	-	0/4/4/4	0/0/0/0
7	GOL	B	805	-	-	0/4/4/4	0/0/0/0
7	GOL	B	807	-	-	0/4/4/4	0/0/0/0
7	GOL	B	808	-	-	0/4/4/4	0/0/0/0
7	GOL	B	809	-	-	0/4/4/4	0/0/0/0

The worst 5 of 38 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	344	HEM	C3B-C4B	-7.75	1.44	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	344	HEM	CAD-C3D	-6.42	1.41	1.54
6	B	344	HEM	C2D-C3D	-4.81	1.40	1.54
6	A	344	HEM	C3B-CAB	-4.70	1.42	1.51
6	B	344	HEM	C4A-CHB	-3.45	1.30	1.39

The worst 5 of 45 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	344	HEM	CAA-CBA-CGA	-6.70	100.46	112.75
6	B	344	HEM	CAA-CBA-CGA	-3.96	105.49	112.75
6	A	344	HEM	CBA-CAA-C2A	-2.81	107.50	112.53
2	A	500	NAG	C2-N2-C7	-2.54	119.77	123.04
2	B	500	NAG	C2-N2-C7	-2.48	119.86	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	600	BMA	1	0
4	A	700	MAN	1	0
3	B	600	BMA	1	0
7	B	805	GOL	1	0
7	B	808	GOL	2	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	339/343 (98%)	-0.27	5 (1%) 76 81	8, 13, 23, 30	0
1	B	338/343 (98%)	-0.22	6 (1%) 71 76	8, 13, 26, 32	0
All	All	677/686 (98%)	-0.24	11 (1%) 74 79	8, 13, 24, 32	0

The worst 5 of 11 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	334	GLY	5.2
1	A	5	GLY	5.1
1	B	334	GLY	4.9
1	B	6	GLY	3.6
1	A	333	SER	3.4

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
7	GOL	B	808	6/6	0.64	0.51	35.35	63,63,63,63	0
7	GOL	B	803	6/6	0.53	0.45	23.01	59,59,59,60	0
7	GOL	B	804	6/6	0.63	0.39	16.93	52,52,52,52	0
7	GOL	B	805	6/6	0.65	0.41	12.17	54,54,55,55	0
7	GOL	B	809	6/6	0.65	0.33	10.68	56,56,56,56	0
7	GOL	A	806	6/6	0.89	0.23	10.63	40,40,40,40	0
7	GOL	A	811	6/6	0.54	0.48	8.85	53,53,54,54	0
7	GOL	B	807	6/6	0.75	0.22	5.99	50,50,50,51	0
7	GOL	A	801	6/6	0.53	0.32	5.42	35,36,37,37	0
7	GOL	A	810	6/6	0.73	0.21	4.41	52,53,53,53	0
3	BMA	B	600	11/12	0.64	0.36	2.82	35,36,37,38	0
4	MAN	A	700	11/12	0.89	0.14	1.71	24,24,24,25	0
4	MAN	B	700	11/12	0.91	0.13	1.56	22,23,24,24	0
3	BMA	A	600	11/12	0.59	0.28	0.90	31,32,33,34	0
2	NAG	B	500	14/15	0.92	0.11	0.27	16,17,18,18	0
5	CA	B	9002	1/1	1.00	0.11	-0.12	11,11,11,11	0
2	NAG	A	500	14/15	0.94	0.08	-0.36	13,13,15,16	0
6	HEM	B	344	43/43	0.98	0.08	-0.38	8,10,12,14	0
6	HEM	A	344	43/43	0.98	0.08	-0.81	8,9,10,12	0
5	CA	A	9001	1/1	1.00	0.09	-0.95	10,10,10,10	0
5	CA	A	9004	1/1	1.00	0.05	-2.27	9,9,9,9	0
5	CA	B	9003	1/1	1.00	0.05	-2.39	8,8,8,8	0
7	GOL	B	800	6/6	0.85	0.13	-	36,36,37,37	0
7	GOL	B	802	6/6	0.71	0.41	-	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.