



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 24, 2022 – 12:13 pm GMT

PDB ID : 5OA5
Title : CELLOBIOHYDROLASE I (CEL7A) FROM HYPOCREA JECORINA
WITH IMPROVED THERMAL STABILITY
Authors : Goedegebuur, F.; Hansson, H.; Karkehabadi, S.; Mikkelsen, N.; Stahlberg, J.;
Sandgren, M.
Deposited on : 2017-06-20
Resolution : 2.10 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.27
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0267
CCP4 : 7.1.010 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.27

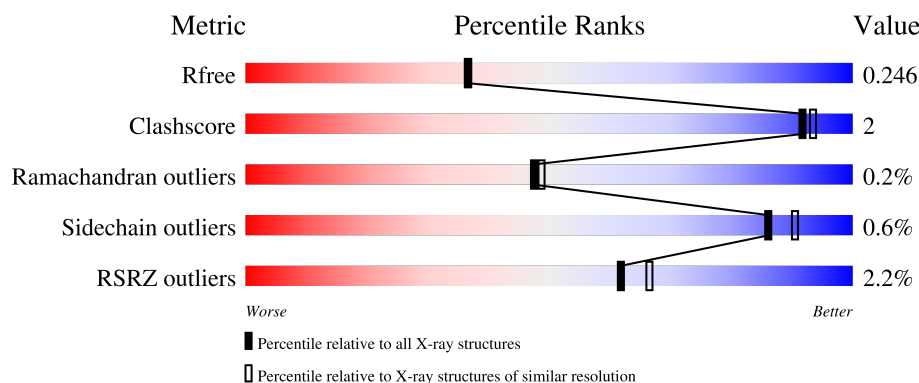
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

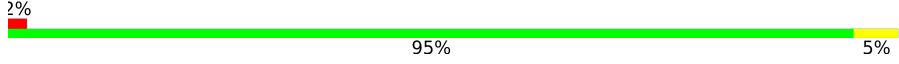
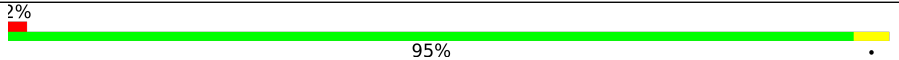
The reported resolution of this entry is 2.10 Å.

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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 $\leq 5\%$. 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	434	
1	B	434	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7133 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 Exoglucanase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	434	Total	C	N	O	S	0	0	0
			3240	2000	534	680	26			
1	B	434	Total	C	N	O	S	0	3	0
			3257	2010	535	686	26			

There are 34 discrepancies between the modelled and reference sequences:

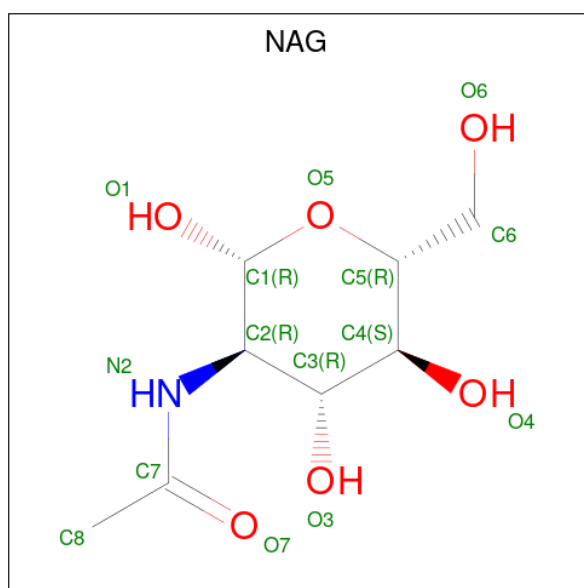
Chain	Residue	Modelled	Actual	Comment	Reference
A	8	PRO	SER	engineered mutation	UNP P62694
A	41	ILE	THR	engineered mutation	UNP P62694
A	49	SER	ASN	engineered mutation	UNP P62694
A	68	THR	ALA	engineered mutation	UNP P62694
A	89	ASP	ASN	engineered mutation	UNP P62694
A	92	THR	SER	engineered mutation	UNP P62694
A	113	ASN	SER	engineered mutation	UNP P62694
A	196	THR	SER	engineered mutation	UNP P62694
A	227	LEU	PRO	engineered mutation	UNP P62694
A	249	LYS	ASP	engineered mutation	UNP P62694
A	255	PRO	THR	engineered mutation	UNP P62694
A	278	PRO	SER	engineered mutation	UNP P62694
A	295	LYS	GLU	engineered mutation	UNP P62694
A	296	PRO	THR	engineered mutation	UNP P62694
A	332	TYR	THR	engineered mutation	UNP P62694
A	403	ASP	VAL	engineered mutation	UNP P62694
A	411	PHE	SER	engineered mutation	UNP P62694
B	8	PRO	SER	engineered mutation	UNP P62694
B	41	ILE	THR	engineered mutation	UNP P62694
B	49	SER	ASN	engineered mutation	UNP P62694
B	68	THR	ALA	engineered mutation	UNP P62694
B	89	ASP	ASN	engineered mutation	UNP P62694
B	92	THR	SER	engineered mutation	UNP P62694
B	113	ASN	SER	engineered mutation	UNP P62694
B	196	THR	SER	engineered mutation	UNP P62694

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Chain	Residue	Modelled	Actual	Comment	Reference
B	227	LEU	PRO	engineered mutation	UNP P62694
B	249	LYS	ASP	engineered mutation	UNP P62694
B	255	PRO	THR	engineered mutation	UNP P62694
B	278	PRO	SER	engineered mutation	UNP P62694
B	295	LYS	GLU	engineered mutation	UNP P62694
B	296	PRO	THR	engineered mutation	UNP P62694
B	332	TYR	THR	engineered mutation	UNP P62694
B	403	ASP	VAL	engineered mutation	UNP P62694
B	411	PHE	SER	engineered mutation	UNP P62694

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
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		
2	B	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 GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	277	Total	O	0	0
			277	277		
4	B	269	Total	O	0	0
			269	269		

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: Exoglucanase 1



- Molecule 1: Exoglucanase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	102.94Å 92.10Å 102.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.16 – 2.10 38.88 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.1 (41.16-2.10) 99.1 (38.88-2.10)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.199 , 0.244 0.200 , 0.246	Depositor DCC
R_{free} test set	2883 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	16.0	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7133	wwPDB-VP
Average B, all atoms (Å ²)	17.0	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 50.84 % 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 6.1087e-05. 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, 0.333 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: PCA, GOL, NAG

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	0.51	0/3313	0.63	0/4517
1	B	0.50	0/3336	0.62	0/4548
All	All	0.50	0/6649	0.63	0/9065

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

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	3240	0	2953	10	0
1	B	3257	0	2967	10	0
2	A	42	0	39	1	0
2	B	42	0	39	0	0
3	B	6	0	8	0	0
4	A	277	0	0	1	0
4	B	269	0	0	1	0
All	All	7133	0	6006	20	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 2.

All (20) 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:247:TYR:O	1:B:248:SER:HB3	2.08	0.54
1:A:227:LEU:HD21	1:A:300:ILE:HD12	1.90	0.53
1:B:181:LYS:HE3	4:B:708:HOH:O	2.09	0.51
1:B:177:PRO:HD2	1:B:208:SER:O	2.11	0.50
1:A:1:PCA:OE	1:A:185:GLY:HA2	2.14	0.48
1:A:68:THR:HB	4:A:635:HOH:O	2.16	0.44
1:A:289:THR:HB	1:A:305:VAL:HB	1.98	0.44
1:B:227:LEU:HD21	1:B:300:ILE:HD12	2.00	0.42
1:A:11:HIS:HB3	1:A:32:VAL:O	2.19	0.42
1:B:63:ASP:HB2	1:B:190:GLU:OE2	2.20	0.41
1:B:289:THR:HB	1:B:305:VAL:HB	2.02	0.41
1:A:17:GLN:HB2	1:A:422:LYS:HG2	2.02	0.41
1:A:186:GLN:O	2:A:501:NAG:H82	2.21	0.41
1:A:27:GLN:HB2	1:A:434:GLY:O	2.20	0.41
1:B:287:LYS:HG2	1:B:288:LEU:N	2.35	0.41
1:A:225:LEU:HD11	1:A:326:LEU:HD13	2.03	0.41
1:A:387:SER:HB3	1:A:393:VAL:HG21	2.02	0.41
1:B:17:GLN:HB2	1:B:422:LYS:HG2	2.03	0.41
1:B:192:TRP:HA	1:B:204:GLY:HA2	2.03	0.41
1:B:139:GLY:HA3	1:B:402:GLY:HA2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	432/434 (100%)	419 (97%)	12 (3%)	1 (0%)	47	49
1	B	435/434 (100%)	425 (98%)	9 (2%)	1 (0%)	47	49
All	All	867/868 (100%)	844 (97%)	21 (2%)	2 (0%)	47	49

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	248	SER
1	A	248	SER

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	356/356 (100%)	353 (99%)	3 (1%)	81	86
1	B	359/356 (101%)	358 (100%)	1 (0%)	92	95
All	All	715/712 (100%)	711 (99%)	4 (1%)	86	90

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	SER
1	A	288	LEU
1	A	398	SER
1	B	288	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	431	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](#)

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	PCA	B	1	1	7,8,9	0.52	0	9,10,12	1.61	1 (11%)
1	PCA	A	1	1	7,8,9	0.48	0	9,10,12	1.86	1 (11%)

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
1	PCA	B	1	1	-	0/0/11/13	0/1/1/1
1	PCA	A	1	1	-	0/0/11/13	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	PCA	CB-CA-C	-4.80	106.09	112.70
1	B	1	PCA	CB-CA-C	-3.96	107.25	112.70

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
1	A	1	PCA	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry

7 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	NAG	A	501	1	14,14,15	0.58	0	17,19,21	1.09	2 (11%)
2	NAG	A	503	1	14,14,15	0.62	0	17,19,21	0.82	0
2	NAG	B	503	1	14,14,15	0.55	0	17,19,21	0.82	0
2	NAG	A	502	1	14,14,15	0.70	0	17,19,21	1.21	1 (5%)
2	NAG	B	501	1	14,14,15	0.53	0	17,19,21	2.30	2 (11%)
2	NAG	B	502	1	14,14,15	0.50	0	17,19,21	1.44	3 (17%)
3	GOL	B	504	-	5,5,5	0.27	0	5,5,5	0.41	0

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	NAG	A	501	1	-	2/6/23/26	0/1/1/1
2	NAG	A	503	1	-	0/6/23/26	0/1/1/1
2	NAG	B	503	1	-	0/6/23/26	0/1/1/1
2	NAG	A	502	1	-	2/6/23/26	0/1/1/1
2	NAG	B	501	1	-	0/6/23/26	0/1/1/1
2	NAG	B	502	1	-	0/6/23/26	0/1/1/1
3	GOL	B	504	-	-	0/4/4/4	-

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NAG	C1-O5-C5	8.54	123.77	112.19
2	B	502	NAG	O5-C5-C6	2.87	111.71	107.20
2	A	502	NAG	O5-C1-C2	-2.66	107.09	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAG	O5-C5-C6	2.49	111.11	107.20
2	B	502	NAG	C3-C4-C5	-2.35	106.05	110.24
2	B	502	NAG	C4-C3-C2	-2.29	107.67	111.02
2	A	501	NAG	C1-O5-C5	2.20	115.17	112.19
2	B	501	NAG	C6-C5-C4	-2.15	107.97	113.00

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NAG	O5-C5-C6-O6
2	A	501	NAG	C4-C5-C6-O6
2	A	502	NAG	C4-C5-C6-O6
2	A	502	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	NAG	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	433/434 (99%)	-0.28	9 (2%) 63 68	7, 15, 32, 73	0
1	B	433/434 (99%)	-0.28	10 (2%) 60 65	7, 15, 36, 64	0
All	All	866/868 (99%)	-0.28	19 (2%) 62 66	7, 15, 32, 73	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	432	PRO	5.9
1	A	247	TYR	4.4
1	B	388	SER	4.4
1	B	247	TYR	4.0
1	A	434	GLY	3.7
1	A	23	GLY	3.5
1	B	434	GLY	3.5
1	A	432	PRO	3.4
1	A	249	LYS	3.2
1	B	384	ASN	3.0
1	B	386	THR	2.9
1	B	249	LYS	2.9
1	B	22	GLY	2.6
1	A	371	TYR	2.6
1	B	387	SER	2.5
1	A	384	ASN	2.5
1	B	101	GLN	2.4
1	A	250	ASN	2.2
1	A	386	THR	2.2

6.2 Non-standard residues in protein, DNA, RNA chains [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, 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	B-factors(Å ²)	Q<0.9
1	PCA	A	1	8/9	0.97	0.07	11,12,12,12	0
1	PCA	B	1	8/9	0.97	0.09	9,10,10,10	0

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, 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	B-factors(Å ²)	Q<0.9
3	GOL	B	504	6/6	0.81	0.17	35,40,41,42	0
2	NAG	B	503	14/15	0.84	0.19	29,32,35,37	0
2	NAG	A	501	14/15	0.85	0.25	32,38,39,41	0
2	NAG	B	501	14/15	0.86	0.26	30,34,37,39	0
2	NAG	B	502	14/15	0.87	0.14	29,34,37,38	0
2	NAG	A	502	14/15	0.91	0.13	21,24,27,27	0
2	NAG	A	503	14/15	0.92	0.16	22,26,30,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.