



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 05:47 pm BST

PDB ID : 2CIO
Title : The high resolution x-ray structure of papain complexed with fragments of the Trypanosoma brucei cysteine protease inhibitor ICP.
Authors : Alphey, M.S.; Hunter, W.N.
Deposited on : 2006-03-24
Resolution : 1.50 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

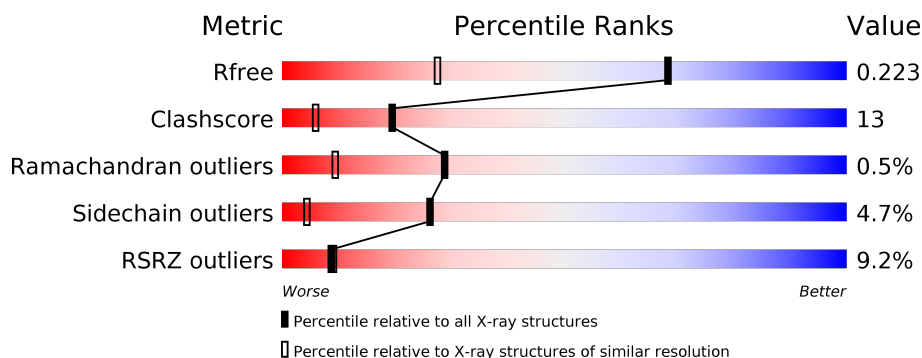
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 1.50 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	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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 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	212	
2	B	121	

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
3	GOL	A	1217	-	-	X	X
4	ACT	A	1218[A]	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 1974 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 PAPAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	17	0
			1753	1122	302	322	7			

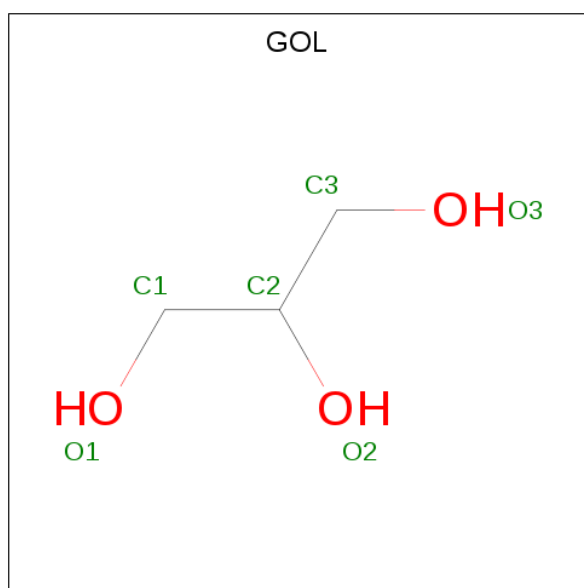
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	GLN	GLU	conflict	UNP P00784

- Molecule 2 is a protein called INHIBITOR OF CYSTEINE PEPTIDASE.

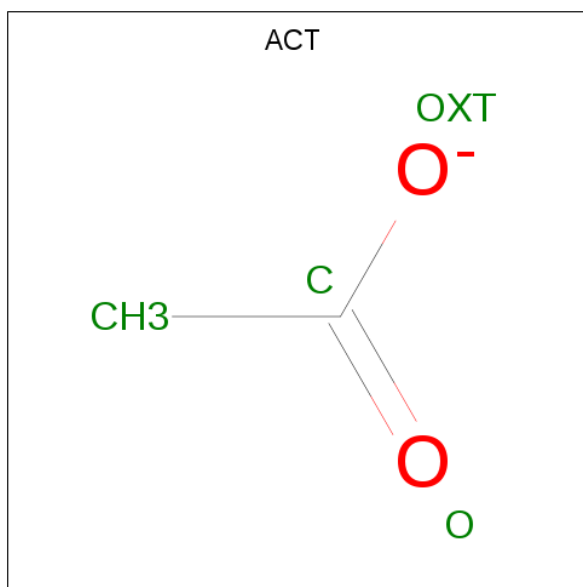
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	7	Total	C	N	O	0	0	2
			32	19	7	6			

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).

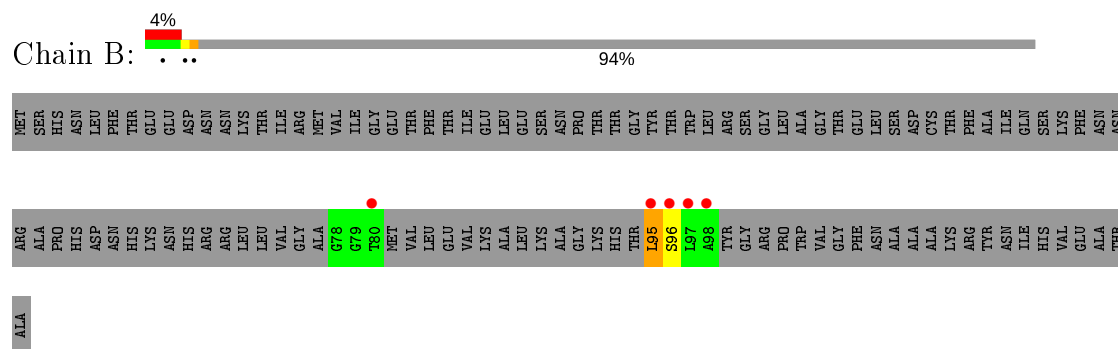


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	1
			8	4	4		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	158	Total	O	0	0
			158	158		
5	B	3	Total	O	0	0
			3	3		

- Molecule 1: PAPAIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	42.32Å 46.12Å 95.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	95.70 – 1.50 46.12 – 1.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (95.70-1.50) 100.0 (46.12-1.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.84 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.177 , 0.225 0.175 , 0.223	Depositor DCC
R_{free} test set	1550 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	15.3	Xtriage
Anisotropy	0.726	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 64.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	1974	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.46% 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: GOL, OCS, ACT

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	0.65	0/1836	0.70	0/2490
2	B	1.33	0/30	1.03	0/38
All	All	0.67	0/1866	0.71	0/2528

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	1753	0	1735	45	0
2	B	32	0	31	3	0
3	A	12	0	16	6	0
4	A	16	0	12	4	0
5	A	158	0	0	14	1
5	B	3	0	0	0	0
All	All	1974	0	1794	46	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 13.

All (46) 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:76:ALA:HB3	5:A:2063:HOH:O	1.25	1.27
1:A:8:ARG:HG2	1:A:13[B]:VAL:HG11	1.32	1.08
1:A:155[B]:ASN:ND2	1:A:197[B]:TYR:CD2	2.26	1.03
1:A:78:TYR:HE1	3:A:1217:GOL:HO2	1.09	0.93
1:A:78:TYR:HE1	3:A:1217:GOL:O2	1.53	0.90
1:A:83:ARG:HH12	4:A:1218[A]:ACT:H1	1.40	0.86
3:A:1217:GOL:H32	5:A:2047:HOH:O	1.78	0.84
1:A:155[A]:ASN:H	1:A:155[A]:ASN:HD22	1.28	0.81
1:A:98[B]:ARG:H	1:A:98[B]:ARG:HH21	1.29	0.78
1:A:109:GLY:HA2	5:A:2099:HOH:O	1.88	0.73
1:A:73[B]:GLN:HA	5:A:2063:HOH:O	1.90	0.71
1:A:156:LYS:O	1:A:156:LYS:HG3	1.90	0.70
1:A:3[B]:GLU:OE1	5:A:2004:HOH:O	2.10	0.70
1:A:41:ARG:CZ	1:A:211:LYS:O	2.40	0.69
1:A:74[A]:LEU:HB2	3:A:1217:GOL:H31	1.76	0.68
1:A:98[B]:ARG:NH2	5:A:2088:HOH:O	2.28	0.66
1:A:76:ALA:CB	5:A:2063:HOH:O	2.04	0.66
1:A:211:LYS:HG2	1:A:212:ASN:HB3	1.77	0.66
1:A:73[A]:GLN:HA	5:A:2063:HOH:O	1.96	0.65
1:A:8:ARG:HG2	1:A:13[B]:VAL:CG1	2.21	0.64
1:A:156:LYS:HD2	5:A:2127:HOH:O	1.99	0.62
1:A:159:HIS:HD1	2:B:95:LEU:N	1.98	0.61
1:A:41:ARG:NH2	1:A:211:LYS:O	2.35	0.60
1:A:109:GLY:CA	5:A:2099:HOH:O	2.49	0.59
1:A:156:LYS:O	1:A:156:LYS:CG	2.49	0.59
1:A:78:TYR:CE1	3:A:1217:GOL:O2	2.39	0.59
1:A:84[B]:ASN:ND2	5:A:2074:HOH:O	2.32	0.59
1:A:59:ARG:NH1	5:A:2050:HOH:O	2.37	0.57
1:A:83:ARG:HH12	4:A:1218[A]:ACT:CH3	2.18	0.55
1:A:98[A]:ARG:NH1	4:A:1220:ACT:OXT	2.42	0.53
1:A:173[A]:ILE:HB	1:A:187:ILE:HG23	1.92	0.52
1:A:74[A]:LEU:HD12	1:A:78:TYR:HB2	1.92	0.52
1:A:146:GLY:HA2	1:A:183:GLU:HG3	1.94	0.50
1:A:108:ASP:OD2	1:A:212:ASN:HA	2.13	0.48
1:A:73[B]:GLN:NE2	5:A:2064:HOH:O	2.48	0.47
1:A:142:GLN:HE22	2:B:95:LEU:HD12	1.79	0.47
1:A:59:ARG:NH2	4:A:1219:ACT:OXT	2.44	0.45
1:A:155[B]:ASN:ND2	1:A:197[B]:TYR:CG	2.81	0.45
1:A:74[A]:LEU:HD13	3:A:1217:GOL:O3	2.17	0.44
1:A:142:GLN:NE2	2:B:95:LEU:HD12	2.32	0.44
1:A:154:GLY:O	1:A:200:CYS:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:LEU:HD13	1:A:71:ALA:HB1	2.00	0.42
1:A:161:VAL:HG21	1:A:173[B]:ILE:HG21	2.02	0.42
1:A:73[A]:GLN:NE2	5:A:2063:HOH:O	2.54	0.41
1:A:69:TRP:O	1:A:73[B]:GLN:HG2	2.21	0.41

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)
5:A:2072:HOH:O	5:A:2140:HOH:O[4_455]	1.92	0.28

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	226/212 (107%)	218 (96%)	7 (3%)	1 (0%)	34	13
2	B	3/121 (2%)	3 (100%)	0	0	100	100
All	All	229/333 (69%)	221 (96%)	7 (3%)	1 (0%)	29	13

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	193	THR

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	186/169 (110%)	179 (96%)	7 (4%)	33 7
2	B	3/99 (3%)	1 (33%)	2 (67%)	0 0
All	All	189/268 (70%)	180 (95%)	9 (5%)	26 4

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	ARG
1	A	77	GLN
1	A	86	TYR
1	A	155[A]	ASN
1	A	155[B]	ASN
1	A	156	LYS
1	A	195	ASN
2	B	95	LEU
2	B	96	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	112	GLN
1	A	195	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	OCS	A	25	1	7,8,9	0.94	0	6,11,13	1.42	1 (16%)

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	OCS	A	25	1	-	2/4/7/9	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	25	OCS	OD1-SG-CB	2.96	110.46	106.94

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	25	OCS	N-CA-CB-SG
1	A	25	OCS	CA-CB-SG-OD2

There are no ring outliers.

No monomer is involved in short contacts.

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	$\# Z > 2$
4	ACT	A	1220	-	1,3,3	1.33	0	0,3,3	0.00	-
4	ACT	A	1219	-	1,3,3	1.25	0	0,3,3	0.00	-
3	GOL	A	1216	-	5,5,5	0.31	0	5,5,5	0.45	0
3	GOL	A	1217	-	5,5,5	0.39	0	5,5,5	0.52	0
4	ACT	A	1218[B]	-	1,3,3	1.47	0	0,3,3	0.00	-
4	ACT	A	1218[A]	-	1,3,3	1.22	0	0,3,3	0.00	-

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	GOL	A	1217	-	-	4/4/4/4	-
3	GOL	A	1216	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1217	GOL	O1-C1-C2-C3
3	A	1217	GOL	C1-C2-C3-O3
3	A	1217	GOL	O2-C2-C3-O3
3	A	1217	GOL	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1220	ACT	1	0
4	A	1219	ACT	1	0
3	A	1217	GOL	6	0
4	A	1218[A]	ACT	2	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	211/212 (99%)	0.25	15 (7%) 16 17	9, 17, 29, 41	2 (0%)
2	B	7/121 (5%)	2.84	5 (71%) 0 0	33, 37, 38, 39	0
All	All	218/333 (65%)	0.33	20 (9%) 9 9	9, 17, 33, 41	2 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	98	ALA	6.5
1	A	197[A]	TYR	5.3
1	A	194	GLY	5.3
1	A	212	ASN	3.7
2	B	97	LEU	3.5
1	A	77	GLN	3.2
1	A	192	GLY	3.0
1	A	98[A]	ARG	2.9
1	A	195	ASN	2.8
1	A	76	ALA	2.6
2	B	95	LEU	2.6
1	A	103	TYR	2.6
1	A	208	TYR	2.6
2	B	96	SER	2.5
2	B	80	THR	2.4
1	A	111	ARG	2.3
1	A	110	VAL	2.2
1	A	74[A]	LEU	2.2
1	A	78	TYR	2.1
1	A	73[A]	GLN	2.0

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

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	OCS	A	25	9/10	0.98	0.06	10,11,16,16	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	A	1217	6/6	0.35	0.64	50,52,52,56	0
3	GOL	A	1216	6/6	0.67	0.14	37,43,45,46	0
4	ACT	A	1219	4/4	0.87	0.16	43,43,43,44	0
4	ACT	A	1220	4/4	0.94	0.09	48,48,48,48	0
4	ACT	A	1218[B]	4/4	0.95	0.12	5,6,6,6	4
4	ACT	A	1218[A]	4/4	0.95	0.12	18,18,20,20	4

6.5 Other polymers ⓘ

There are no such residues in this entry.