



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:48 AM UTC

PDB ID : 3BJE / pdb_00003bje
Title : Crystal structure of Trypanosoma brucei nucleoside phosphorylase shows uridine phosphorylase activity
Authors : Larson, E.T.; Merritt, E.A.; Structural Genomics of Pathogenic Protozoa Consortium (SGPP)
Deposited on : 2007-12-03
Resolution : 1.44 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

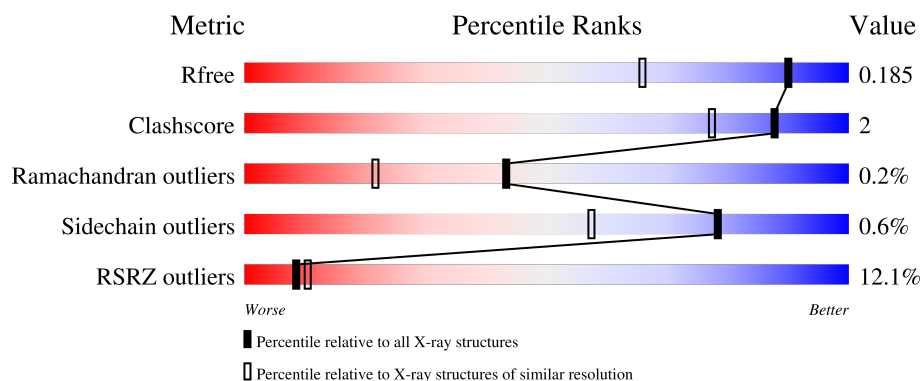
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.44 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}	180053	3234 (1.46-1.42)
Clashscore	190562	3289 (1.46-1.42)
Ramachandran outliers	187476	3248 (1.46-1.42)
Sidechain outliers	187428	3248 (1.46-1.42)
RSRZ outliers	180081	3234 (1.46-1.42)

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	349	12% 91% 6%
1	B	349	10% 88% 5% 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5601 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.

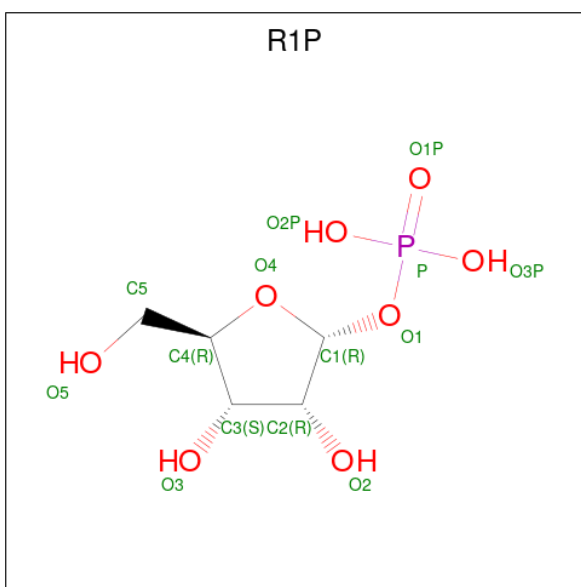
- Molecule 1 is a protein called Nucleoside phosphorylase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	327	Total	C	N	O	S	0	5	0
			2521	1568	450	484	19			
1	B	327	Total	C	N	O	S	0	8	0
			2526	1573	444	489	20			

There are 16 discrepancies between the modelled and reference sequences:

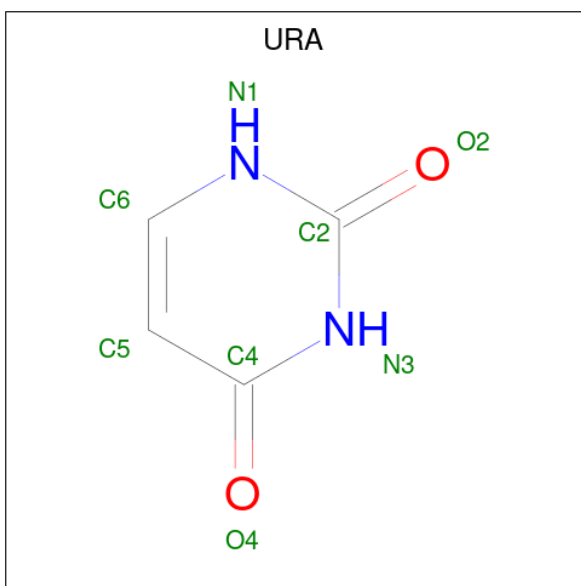
Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	expression tag	UNP Q57VZ2
A	-6	ALA	-	expression tag	UNP Q57VZ2
A	-5	HIS	-	expression tag	UNP Q57VZ2
A	-4	HIS	-	expression tag	UNP Q57VZ2
A	-3	HIS	-	expression tag	UNP Q57VZ2
A	-2	HIS	-	expression tag	UNP Q57VZ2
A	-1	HIS	-	expression tag	UNP Q57VZ2
A	0	HIS	-	expression tag	UNP Q57VZ2
B	-7	MET	-	expression tag	UNP Q57VZ2
B	-6	ALA	-	expression tag	UNP Q57VZ2
B	-5	HIS	-	expression tag	UNP Q57VZ2
B	-4	HIS	-	expression tag	UNP Q57VZ2
B	-3	HIS	-	expression tag	UNP Q57VZ2
B	-2	HIS	-	expression tag	UNP Q57VZ2
B	-1	HIS	-	expression tag	UNP Q57VZ2
B	0	HIS	-	expression tag	UNP Q57VZ2

- Molecule 2 is 1-O-phosphono-alpha-D-ribofuranose (CCD ID: R1P) (formula: C₅H₁₁O₈P).



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

- Molecule 3 is URACIL (CCD ID: URA) (formula: $C_4H_4N_2O_2$).



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

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Ca 1	0	0

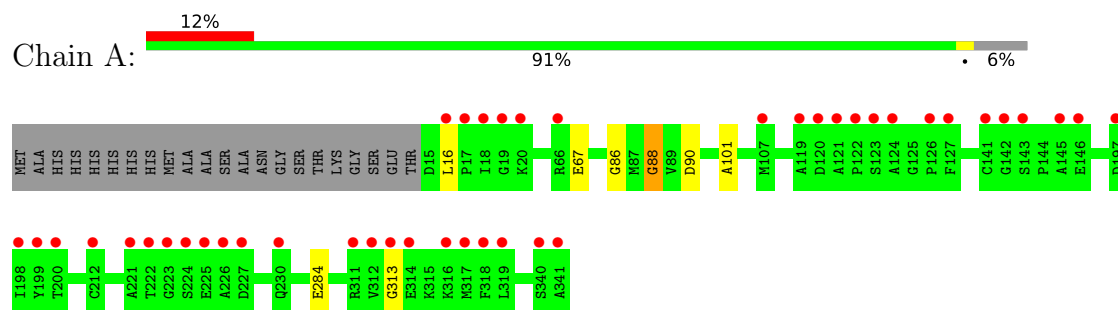
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	234	Total 238	O 238	0	4
5	B	267	Total 271	O 271	0	4

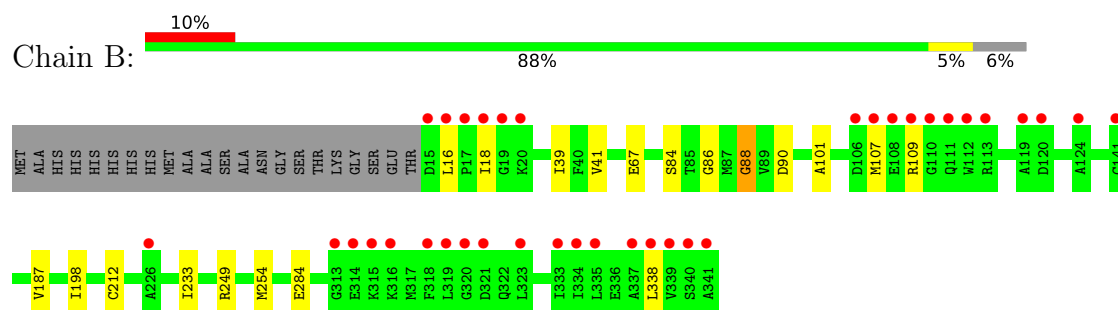
3 Residue-property plots [i](#)

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: Nucleoside phosphorylase, putative



- Molecule 1: Nucleoside phosphorylase, putative



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.02Å 95.39Å 63.48Å 90.00° 105.91° 90.00°	Depositor
Resolution (Å)	35.38 – 1.44 35.38 – 1.44	Depositor EDS
% Data completeness (in resolution range)	94.7 (35.38-1.44) 94.7 (35.38-1.44)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 1.44Å)	Xtriage
Refinement program	REFMAC refmac_5.2.0019	Depositor
R, R_{free}	0.155 , 0.184 0.158 , 0.185	Depositor DCC
R_{free} test set	6072 reflections (4.67%)	wwPDB-VP
Wilson B-factor (Å ²)	14.8	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5601	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.87% 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: CA, URA, R1P

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	0.59	0/2572	0.76	1/3468 (0.0%)
1	B	0.61	0/2589	0.77	1/3492 (0.0%)
All	All	0.60	0/5161	0.76	2/6960 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	GLY	N-CA-C	8.24	119.78	111.95
1	B	88	GLY	N-CA-C	5.67	117.34	111.95

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	2521	0	2537	7	0
1	B	2526	0	2542	16	0
2	A	14	0	0	0	0
2	B	14	0	0	0	0
3	A	8	0	3	0	0
3	B	8	0	3	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	238	0	0	0	0
5	B	271	0	0	1	0
All	All	5601	0	5085	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:16:LEU:HD21	1:B:101:ALA:CB	2.14	0.77
1:A:67:GLU:HG3	1:A:86:GLY:HA3	1.74	0.69
1:B:16:LEU:HD21	1:B:101:ALA:HB1	1.73	0.69
1:B:67:GLU:HG3	1:B:86:GLY:HA3	1.80	0.63
1:A:67:GLU:HB3	1:B:67:GLU:HB3	1.81	0.62
1:B:16:LEU:HD21	1:B:101:ALA:HB2	1.82	0.61
1:B:254[B]:MET:HE3	5:B:739:HOH:O	2.01	0.60
1:A:16:LEU:HD21	1:A:101:ALA:CB	2.32	0.59
1:B:187:VAL:HG11	1:B:198:ILE:CD1	2.34	0.58
1:A:16:LEU:HD21	1:A:101:ALA:HB2	1.87	0.56
1:B:18:ILE:HD13	1:B:107:MET:HE1	1.91	0.52
1:B:187:VAL:HG11	1:B:198:ILE:HD13	1.93	0.51
1:B:212:CYS:SG	1:B:233:ILE:HD11	2.57	0.45
1:B:249:ARG:HB3	1:B:254[B]:MET:HE2	2.00	0.44
1:A:90:ASP:OD2	1:B:88:GLY:HA3	2.19	0.43
1:A:16:LEU:HD21	1:A:101:ALA:HB1	2.00	0.42
1:A:88:GLY:HA3	1:B:90:ASP:OD2	2.20	0.41
1:B:39:ILE:HG13	1:B:338:LEU:HD22	2.03	0.41
1:B:41:VAL:O	1:B:84:SER:HA	2.21	0.41
1:B:187:VAL:CG1	1:B:198:ILE:CD1	3.00	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	330/349 (95%)	322 (98%)	7 (2%)	1 (0%)	36	17
1	B	333/349 (95%)	325 (98%)	8 (2%)	0	100	100
All	All	663/698 (95%)	647 (98%)	15 (2%)	1 (0%)	43	22

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	313	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	273/284 (96%)	272 (100%)	1 (0%)	84	69
1	B	276/284 (97%)	274 (99%)	2 (1%)	76	53
All	All	549/568 (97%)	546 (100%)	3 (0%)	78	63

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	284	GLU
1	B	109	ARG
1	B	284	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	97	ASN
1	A	180	GLN
1	A	230	GLN
1	B	97	ASN

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Mol	Chain	Res	Type
1	B	181	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	$\# Z > 2$
2	R1P	A	401	-	13,14,14	0.85	0	21,21,21	1.71	4 (19%)
3	URA	A	501	-	8,8,8	1.84	4 (50%)	10,10,10	2.58	3 (30%)
2	R1P	B	401	-	13,14,14	0.92	0	21,21,21	1.89	4 (19%)
3	URA	B	501	-	8,8,8	1.70	2 (25%)	10,10,10	2.75	5 (50%)

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	R1P	A	401	-	-	0/6/23/23	0/1/1/1
3	URA	A	501	-	-	-	0/1/1/1
2	R1P	B	401	-	-	0/6/23/23	0/1/1/1
3	URA	B	501	-	-	-	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	URA	C2-N3	-2.82	1.32	1.37
3	B	501	URA	C4-N3	-2.70	1.34	1.38
3	A	501	URA	C5-C4	-2.37	1.38	1.43
3	B	501	URA	C6-N1	-2.12	1.32	1.36
3	A	501	URA	C4-N3	-2.04	1.35	1.38
3	A	501	URA	O2-C2	2.02	1.27	1.23

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	R1P	O1-C1-C2	6.13	116.60	106.65
2	A	401	R1P	O1-C1-C2	5.12	114.95	106.65
3	B	501	URA	N1-C2-N3	4.91	120.35	115.17
3	B	501	URA	C4-N3-C2	-4.83	120.87	125.55
3	A	501	URA	N1-C2-N3	4.35	119.76	115.17
3	A	501	URA	C4-N3-C2	-4.34	121.35	125.55
3	A	501	URA	C5-C4-N3	3.60	119.84	114.80
3	B	501	URA	O2-C2-N1	-3.15	119.54	122.79
3	B	501	URA	C5-C4-N3	2.92	118.88	114.80
2	B	401	R1P	O1-P-O1P	-2.87	99.10	109.33
2	A	401	R1P	O1-P-O1P	-2.68	99.77	109.33
2	B	401	R1P	O4-C1-O1	2.36	113.66	109.16
2	A	401	R1P	O4-C1-O1	2.34	113.63	109.16
3	B	501	URA	C6-N1-C2	-2.32	120.98	122.40
2	B	401	R1P	O4-C1-C2	-2.24	102.12	104.98
2	A	401	R1P	O3P-P-O2P	2.10	115.67	107.80

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

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	327/349 (93%)	0.97	43 (13%) 7 9	10, 20, 35, 51	5 (1%)
1	B	327/349 (93%)	0.75	36 (11%) 10 12	10, 17, 33, 48	8 (2%)
All	All	654/698 (93%)	0.86	79 (12%) 8 10	10, 19, 34, 51	13 (1%)

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	112	TRP	8.4
1	B	16	LEU	6.0
1	B	18	ILE	5.4
1	A	312	VAL	5.0
1	B	338	LEU	5.0
1	B	113	ARG	4.9
1	B	341	ALA	4.8
1	B	339	VAL	4.3
1	B	319	LEU	4.2
1	B	337	ALA	4.0
1	A	318	PHE	3.9
1	B	19	GLY	3.8
1	A	311	ARG	3.8
1	B	109	ARG	3.8
1	B	17	PRO	3.8
1	A	198	ILE	3.7
1	A	199	TYR	3.7
1	B	334	ILE	3.6
1	A	221	ALA	3.6
1	B	110	GLY	3.6
1	A	319	LEU	3.5
1	B	320	GLY	3.5
1	A	224	SER	3.4
1	B	111	GLN	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	16	LEU	3.4
1	B	315	LYS	3.3
1	A	313	GLY	3.2
1	B	107	MET	3.2
1	A	200	THR	3.1
1	B	314	GLU	3.1
1	B	335	LEU	3.1
1	B	119	ALA	3.1
1	A	18	ILE	3.0
1	A	226	ALA	3.0
1	B	106	ASP	3.0
1	B	318	PHE	3.0
1	A	316	LYS	2.9
1	A	19	GLY	2.8
1	A	227	ASP	2.8
1	A	141	CYS	2.8
1	B	124	ALA	2.8
1	A	20	LYS	2.8
1	B	340[A]	SER	2.8
1	B	313	GLY	2.7
1	B	141	CYS	2.7
1	A	124	ALA	2.7
1	B	15	ASP	2.7
1	A	197	ASP	2.6
1	A	120	ASP	2.5
1	A	122	PRO	2.5
1	A	121	ALA	2.5
1	A	317	MET	2.5
1	B	108	GLU	2.5
1	A	66	ARG	2.5
1	A	145	ALA	2.4
1	A	341	ALA	2.4
1	A	314	GLU	2.4
1	A	146	GLU	2.4
1	A	127	PHE	2.4
1	A	123	SER	2.4
1	A	126	PRO	2.3
1	B	120	ASP	2.3
1	B	316	LYS	2.3
1	A	142	GLY	2.3
1	A	119	ALA	2.2
1	B	321	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	230	GLN	2.2
1	A	107	MET	2.2
1	A	223	GLY	2.1
1	A	143	SER	2.1
1	B	20	LYS	2.1
1	A	225	GLU	2.1
1	A	17	PRO	2.1
1	B	323	LEU	2.1
1	A	212	CYS	2.1
1	B	226	ALA	2.0
1	B	333	ILE	2.0
1	A	340	SER	2.0
1	A	222	THR	2.0

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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	R1P	A	401	14/14	0.94	0.13	19,24,27,28	0
2	R1P	B	401	14/14	0.95	0.14	19,24,28,31	0
3	URA	A	501	8/8	0.96	0.12	16,19,22,23	0
3	URA	B	501	8/8	0.97	0.12	14,17,20,20	0
4	CA	B	402	1/1	1.00	0.15	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.