



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 10, 2020 – 12:10 PM JST

PDB ID : 7C1X  
Title : Unliganded structure of Pseudouridine kinase (PUKI) from Arabidopsis thaliana  
Authors : Kim, S.H.; Rhee, S.  
Deposited on : 2020-05-06  
Resolution : 2.39 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](mailto: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.

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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 : 2.14.6  
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) : 2.14.6

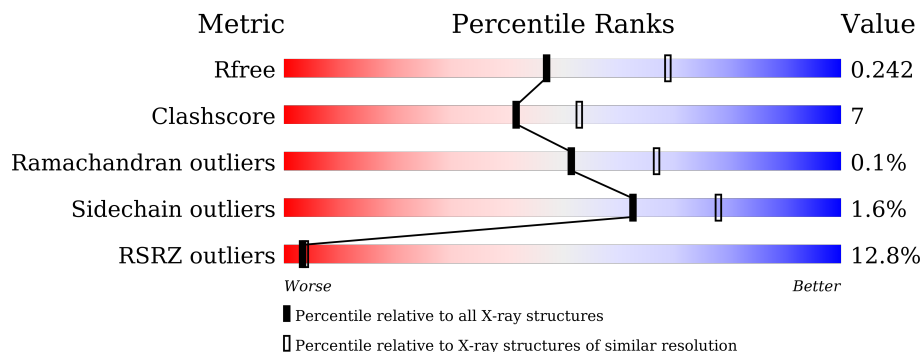
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.39 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	5509 (2.40-2.36)
Clashscore	141614	6082 (2.40-2.36)
Ramachandran outliers	138981	5973 (2.40-2.36)
Sidechain outliers	138945	5975 (2.40-2.36)
RSRZ outliers	127900	5397 (2.40-2.36)

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  $\geq 3$ , 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	378	 10% 78% 16% 7%
1	B	378	 13% 76% 16% 7%

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5225 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 PfkB-like carbohydrate kinase family protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	353	2599	1670	425	488	6	10	0	0	0
1	B	353	2597	1666	429	487	6	9	0	0	0

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Na	0	0
			1	1		
2	A	1	Total	Na	0	0
			1	1		

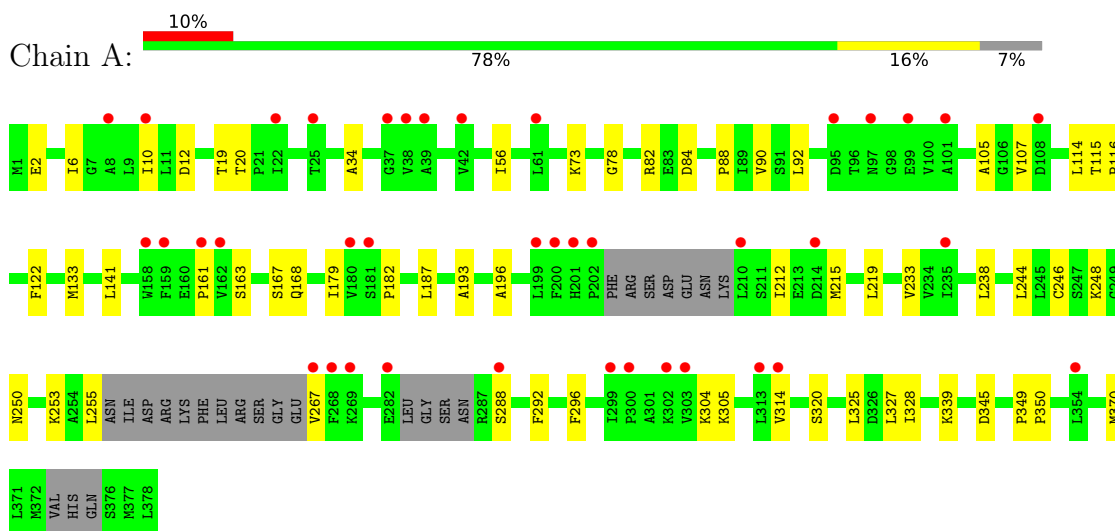
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	10	Total	O	0	0
			10	10		
3	B	17	Total	O	0	0
			17	17		

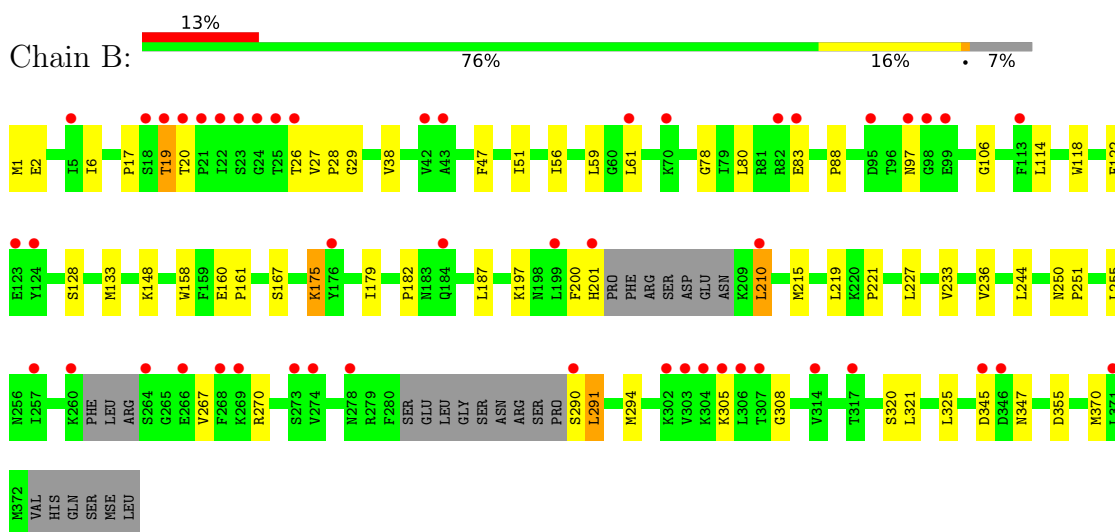
### 3 Residue-property plots

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: PfkB-like carbohydrate kinase family protein



- Molecule 1: PfkB-like carbohydrate kinase family protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	80.69Å 48.35Å 92.75Å 90.00° 109.14° 90.00°	Depositor
Resolution (Å)	29.93 – 2.39 29.93 – 2.39	Depositor EDS
% Data completeness (in resolution range)	98.2 (29.93-2.39) 88.6 (29.93-2.39)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.83 (at 2.39Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.208 , 0.242 0.211 , 0.242	Depositor DCC
$R_{free}$ test set	2000 reflections (7.47%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	44.1	Xtrriage
Anisotropy	0.463	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 54.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5225	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: NA

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	0.23	0/2632	0.45	0/3561
1	B	0.25	0/2630	0.50	0/3560
All	All	0.24	0/5262	0.48	0/7121

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	2599	0	2689	38	1
1	B	2597	0	2681	38	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	10	0	0	0	0
3	B	17	0	0	0	0
All	All	5225	0	5370	74	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 7.

All (74) 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:305:LYS:HE2	1:B:345:ASP:OD1	1.88	0.74
1:A:133:MSE:HE3	1:A:314:VAL:HG11	1.76	0.68
1:B:179:ILE:HG12	1:B:233:VAL:HB	1.75	0.68
1:B:1:MSE:N	1:B:128:SER:O	2.29	0.65
1:B:290:SER:OG	1:B:291:LEU:N	2.29	0.64
1:A:253:LYS:NZ	1:A:288:SER:OG	2.34	0.61
1:B:270:ARG:NH1	1:B:355:ASP:OD1	2.36	0.58
1:A:92:LEU:HD21	1:B:26:THR:HG22	1.85	0.58
1:A:320:SER:HB3	1:A:325:LEU:HD12	1.87	0.57
1:A:10:ILE:HG21	1:A:90:VAL:HG23	1.86	0.57
1:B:175:LYS:H	1:B:175:LYS:HD3	1.69	0.56
1:A:105:ALA:O	1:B:28:PRO:HA	2.06	0.56
1:B:320:SER:HB3	1:B:325:LEU:HD12	1.88	0.55
1:B:270:ARG:HH11	1:B:355:ASP:HA	1.71	0.55
1:A:187:LEU:HD23	1:A:219:LEU:HD22	1.90	0.53
1:B:227:LEU:HD11	1:B:291:LEU:HD23	1.90	0.53
1:B:133:MSE:HG2	1:B:158:TRP:HB3	1.91	0.51
1:B:187:LEU:HD23	1:B:219:LEU:HD22	1.94	0.51
1:B:6:ILE:HG12	1:B:56:ILE:HB	1.93	0.50
1:B:19:THR:OG1	1:B:20:THR:N	2.44	0.50
1:A:90:VAL:HG22	1:A:105:ALA:HB2	1.94	0.50
1:A:179:ILE:HG12	1:A:233:VAL:HB	1.93	0.50
1:B:20:THR:O	1:B:27:VAL:HG21	2.12	0.50
1:B:200:PHE:HE1	1:B:221:PRO:HB2	1.77	0.49
1:B:215:MSE:HE3	1:B:219:LEU:HG	1.93	0.49
1:A:82:ARG:NH1	1:A:84:ASP:OD2	2.46	0.49
1:A:161:PRO:HB3	1:A:167:SER:HB2	1.95	0.48
1:B:161:PRO:HG2	1:B:182:PRO:HB3	1.96	0.48
1:A:246:CYS:HA	1:A:292:PHE:O	2.15	0.47
1:B:17:PRO:HA	1:B:29:GLY:HA3	1.96	0.47
1:B:255:LEU:HD13	1:B:321:LEU:HB3	1.97	0.47
1:A:163:SER:O	1:A:167:SER:HB3	2.15	0.47
1:A:19:THR:OG1	1:A:20:THR:N	2.47	0.47
1:B:210:LEU:HD22	1:B:215:MSE:SE	2.65	0.47
1:B:175:LYS:HD3	1:B:175:LYS:N	2.29	0.46
1:A:233:VAL:HG11	1:A:327:LEU:HD21	1.97	0.46
1:A:179:ILE:HD11	1:A:255:LEU:HD11	1.96	0.46
1:A:12:ASP:HB2	1:A:34:ALA:O	2.15	0.46
1:A:296:PHE:HZ	1:A:328:ILE:HG23	1.80	0.46
1:A:84:ASP:N	1:A:84:ASP:OD1	2.46	0.46
1:B:78:GLY:HA2	1:B:122:PHE:CZ	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:GLY:HA2	1:A:122:PHE:CZ	2.51	0.45
1:A:305:LYS:NZ	1:A:345:ASP:O	2.49	0.45
1:A:304:LYS:HA	1:A:304:LYS:HD2	1.70	0.45
1:B:161:PRO:HB3	1:B:167:SER:HB2	1.99	0.45
1:A:238:LEU:HD11	1:A:244:LEU:HD13	1.99	0.45
1:A:114:LEU:HD12	1:A:114:LEU:HA	1.77	0.45
1:A:292:PHE:HE1	1:A:370:MSE:HE2	1.82	0.44
1:B:114:LEU:HA	1:B:114:LEU:HD12	1.78	0.44
1:B:294:MSE:HG2	1:B:370:MSE:SE	2.67	0.44
1:A:88:PRO:HB3	1:A:107:VAL:HG12	1.99	0.44
1:A:2:GLU:H	1:A:2:GLU:HG2	1.57	0.44
1:A:168:GLN:HG2	1:A:193:ALA:HB2	2.00	0.44
1:B:236:VAL:HB	1:B:244:LEU:HB2	2.00	0.43
1:A:248:LYS:HE2	1:A:248:LYS:HB3	1.73	0.43
1:A:161:PRO:HG2	1:A:182:PRO:HB3	2.01	0.43
1:B:38:VAL:HG11	1:B:160:GLU:HG2	2.01	0.43
1:B:61:LEU:HD12	1:B:83:GLU:HG2	2.01	0.43
1:B:80:LEU:HB2	1:B:118:TRP:CE2	2.55	0.42
1:B:97:ASN:OD1	1:B:97:ASN:N	2.52	0.42
1:A:6:ILE:HG12	1:A:56:ILE:HB	2.01	0.42
1:A:250:ASN:HB3	1:A:253:LYS:HG3	2.02	0.42
1:B:88:PRO:HA	1:B:106:GLY:O	2.20	0.42
1:B:2:GLU:HG2	1:B:2:GLU:H	1.60	0.42
1:B:250:ASN:HA	1:B:251:PRO:HD3	1.81	0.42
1:A:116:PRO:HG2	1:A:141:LEU:HG	2.02	0.41
1:A:73:LYS:HD3	1:A:73:LYS:HA	1.78	0.41
1:A:292:PHE:CE1	1:A:370:MSE:HE2	2.56	0.41
1:B:308:GLY:HA3	1:B:347:ASN:O	2.21	0.41
1:B:59:LEU:HD23	1:B:59:LEU:HA	1.91	0.41
1:A:161:PRO:O	1:A:163:SER:N	2.49	0.41
1:A:212:ILE:HD13	1:A:215:MSE:HE3	2.02	0.40
1:A:349:PRO:HA	1:A:350:PRO:HD3	1.92	0.40
1:B:47:PHE:HA	1:B:51:ILE:O	2.22	0.40

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)
1:A:115:THR:OG1	1:A:196:ALA:O[2_847]	2.15	0.05



## 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	343/378 (91%)	335 (98%)	8 (2%)	0	100	100
1	B	345/378 (91%)	335 (97%)	9 (3%)	1 (0%)	41	53
All	All	688/756 (91%)	670 (97%)	17 (2%)	1 (0%)	51	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	19	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	287/304 (94%)	285 (99%)	2 (1%)	84	92
1	B	284/304 (93%)	277 (98%)	7 (2%)	47	65
All	All	571/608 (94%)	562 (98%)	9 (2%)	62	78

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

Mol	Chain	Res	Type
1	A	267	VAL
1	A	339	LYS
1	B	148	LYS
1	B	175	LYS
1	B	197	LYS

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Mol	Chain	Res	Type
1	B	201	HIS
1	B	210	LEU
1	B	267	VAL
1	B	291	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	343/378 (90%)	0.70	39 (11%) <b>5</b>   <b>5</b>	34, 50, 76, 92	0
1	B	344/378 (91%)	0.86	49 (14%) <b>2</b>   <b>3</b>	40, 57, 90, 102	0
All	All	687/756 (90%)	0.78	88 (12%) <b>3</b>   <b>4</b>	34, 54, 83, 102	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	23	SER	6.3
1	B	304	LYS	6.3
1	B	303	VAL	6.0
1	B	305	LYS	5.0
1	B	176	TYR	4.7
1	B	345	ASP	4.3
1	B	306	LEU	4.1
1	B	19	THR	4.0
1	A	267	VAL	4.0
1	B	201	HIS	4.0
1	B	22	ILE	4.0
1	B	25	THR	3.8
1	B	20	THR	3.8
1	A	61	LEU	3.8
1	A	22	ILE	3.6
1	B	24	GLY	3.6
1	A	302	LYS	3.5
1	A	210	LEU	3.5
1	B	260	LYS	3.5
1	B	199	LEU	3.4
1	B	97	ASN	3.4
1	A	97	ASN	3.4
1	B	269	LYS	3.3
1	A	354	LEU	3.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	B	266	GLU	3.2
1	A	201	HIS	3.2
1	A	200	PHE	3.2
1	B	98	GLY	3.1
1	A	214	ASP	3.0
1	A	314	VAL	2.9
1	B	95	ASP	2.9
1	A	42	VAL	2.9
1	B	70	LYS	2.9
1	B	273	SER	2.9
1	B	274	VAL	2.9
1	A	181	SER	2.8
1	A	159	PHE	2.8
1	B	82	ARG	2.8
1	B	290	SER	2.8
1	A	199	LEU	2.8
1	B	257	ILE	2.8
1	A	313	LEU	2.7
1	A	300	PRO	2.7
1	B	21	PRO	2.7
1	B	26	THR	2.7
1	B	184	GLN	2.7
1	B	210	LEU	2.7
1	B	314	VAL	2.7
1	A	288	SER	2.6
1	A	101	ALA	2.6
1	A	39	ALA	2.6
1	A	269	LYS	2.6
1	A	268	PHE	2.6
1	B	124	TYR	2.6
1	A	161	PRO	2.6
1	B	268	PHE	2.6
1	B	302	LYS	2.6
1	B	83	GLU	2.5
1	B	18	SER	2.5
1	A	38	VAL	2.5
1	A	99	GLU	2.4
1	B	99	GLU	2.4
1	B	307	THR	2.4
1	B	264	SER	2.4
1	A	108	ASP	2.4
1	A	158	TRP	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	8	ALA	2.4
1	A	299	ILE	2.3
1	A	95	ASP	2.3
1	B	61	LEU	2.3
1	B	346	ASP	2.2
1	A	303	VAL	2.2
1	A	37	GLY	2.2
1	B	278	ASN	2.2
1	A	282	GLU	2.2
1	B	113	PHE	2.2
1	B	123	GLU	2.1
1	B	5	ILE	2.1
1	A	202	PRO	2.1
1	A	162	VAL	2.1
1	B	43	ALA	2.1
1	B	371	LEU	2.1
1	A	10	ILE	2.1
1	A	235	ILE	2.1
1	B	42	VAL	2.0
1	A	25	THR	2.0
1	B	317	THR	2.0
1	A	180	VAL	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 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	NA	B	401	1/1	0.91	0.14	66,66,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NA	A	401	1/1	0.95	0.07	67,67,67,67	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.