



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 21, 2022 – 12:23 pm GMT

PDB ID : 7QLJ  
Title : Trans structure of rsKiiro Illuminated at 290 K  
Authors : van Thor, J.J.; Baxter, J.M.  
Deposited on : 2021-12-20  
Resolution : 1.02 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.

The types of validation reports are described at

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

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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 : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.31.3  
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) : 2.31.3

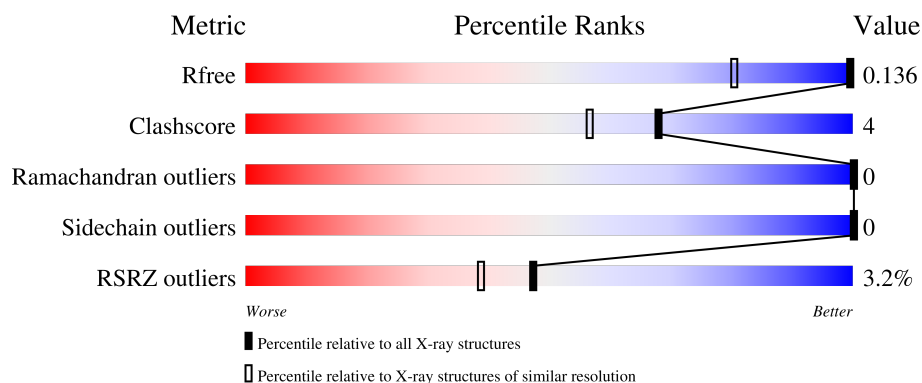
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

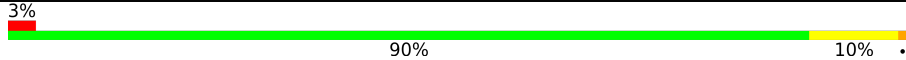
The reported resolution of this entry is 1.02 Å.

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	1188 (1.08-0.96)
Clashscore	141614	1253 (1.08-0.96)
Ramachandran outliers	138981	1178 (1.08-0.96)
Sidechain outliers	138945	1180 (1.08-0.96)
RSRZ outliers	127900	1158 (1.08-0.96)

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  $\geq 3$ , 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	220	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4194 atoms, of which 1951 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.

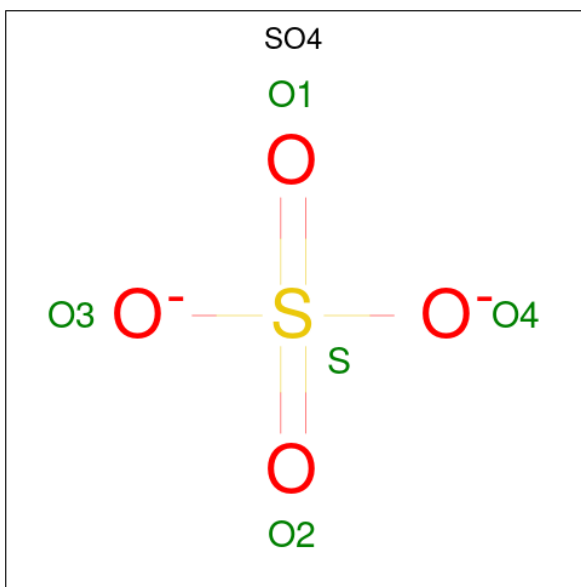
- Molecule 1 is a protein called rsKiir.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	220	3929	1270	1951	334	362	12	65	30	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	LYS	ASN	engineered mutation	UNP Q5S6Z9
A	64	PIA	HIS	chromophore	UNP Q5S6Z9
A	64	PIA	TYR	chromophore	UNP Q5S6Z9
A	64	PIA	GLY	chromophore	UNP Q5S6Z9
A	70	LYS	GLU	engineered mutation	UNP Q5S6Z9
A	74	ASN	HIS	engineered mutation	UNP Q5S6Z9
A	102	ASN	ILE	engineered mutation	UNP Q5S6Z9
A	121	TYR	HIS	engineered mutation	UNP Q5S6Z9
A	123	THR	VAL	engineered mutation	UNP Q5S6Z9
A	157	VAL	ILE	engineered mutation	UNP Q5S6Z9
A	158	GLU	THR	engineered mutation	UNP Q5S6Z9
A	159	THR	MET	engineered mutation	UNP Q5S6Z9
A	189	ALA	TYR	engineered mutation	UNP Q5S6Z9
A	222	ASN	-	expression tag	UNP Q5S6Z9

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

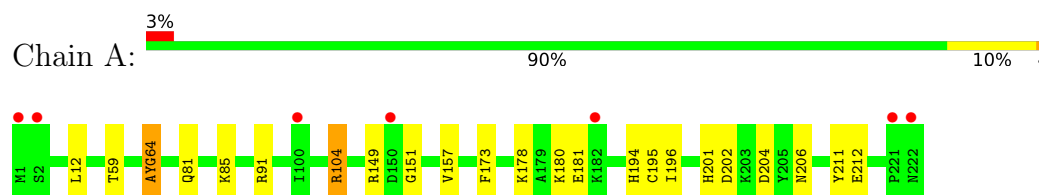
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	260	Total	O	0	3
			260	260		

### 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: rsKiir



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	38.75Å 73.52Å 78.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.05 – 1.02 39.05 – 1.02	Depositor EDS
% Data completeness (in resolution range)	98.9 (39.05-1.02) 98.9 (39.05-1.02)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.81 (at 1.02Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.119 , 0.137 0.119 , 0.136	Depositor DCC
$R_{free}$ test set	5501 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	7.8	Xtriage
Anisotropy	0.113	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.44 , 47.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	4194	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.30% 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: PIA, SO4

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	0.89	4/2038 (0.2%)	1.01	6/2739 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	212	GLU	CD-OE1	-7.91	1.17	1.25
1	A	204	ASP	CG-OD2	-5.70	1.12	1.25
1	A	151	GLY	CA-C	5.26	1.60	1.51
1	A	181	GLU	CD-OE1	5.19	1.31	1.25

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	ARG	NE-CZ-NH2	-10.84	114.88	120.30
1	A	149	ARG	NH1-CZ-NH2	7.34	127.48	119.40
1	A	104	ARG	NE-CZ-NH2	-7.26	116.67	120.30
1	A	211	TYR	CB-CG-CD1	5.87	124.52	121.00
1	A	104	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	A	149	ARG	NE-CZ-NH1	-5.40	117.60	120.30

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	1978	1951	1943	15	0
2	A	5	0	0	0	0
3	A	260	0	0	4	0
All	All	2243	1951	1943	15	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 4.

All (15) 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:104:ARG:NH2	3:A:402:HOH:O	2.31	0.62
1:A:64[C]:PIA:OH	1:A:157:VAL:HG21	2.01	0.61
1:A:195[B]:CYS:SG	3:A:491:HOH:O	2.32	0.58
1:A:64[C]:PIA:OH	1:A:157:VAL:CG2	2.53	0.56
1:A:81[B]:GLN:NE2	3:A:404:HOH:O	2.38	0.55
1:A:59:THR:HB	1:A:64[A]:PIA:HD1	1.91	0.51
1:A:178[B]:LYS:HG3	3:A:501:HOH:O	2.08	0.51
1:A:91:ARG:HD3	1:A:173[A]:PHE:HD2	1.75	0.50
1:A:202:ASP:OD2	1:A:206[B]:ASN:HB3	2.17	0.45
1:A:85:LYS:O	1:A:180[B]:LYS:HD2	2.18	0.43
1:A:59:THR:HB	1:A:64[A]:PIA:CD1	2.48	0.43
1:A:201:HIS:HD2	1:A:202:ASP:O	2.02	0.42
1:A:194[B]:HIS:CE1	1:A:196:ILE:HB	2.55	0.41
1:A:157:VAL:HB	1:A:173[A]:PHE:HB2	2.02	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	244/220 (111%)	243 (100%)	1 (0%)	0	<b>100</b> <b>100</b>



There are no Ramachandran outliers to report.

### 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	216/187 (116%)	216 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	201	HIS

### 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	PIA	A	64[C]	1	21,21,22	1.65	6 (28%)	27,29,31	2.64	10 (37%)
1	PIA	A	64[A]	1	21,21,22	1.66	3 (14%)	27,29,31	3.40	15 (55%)

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	PIA	A	64[C]	1	-	3/8/27/28	0/2/2/2
1	PIA	A	64[A]	1	-	6/8/27/28	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	64[A]	PIA	CA2-C2	4.97	1.53	1.48
1	A	64[C]	PIA	CA1-C1	-3.04	1.47	1.51
1	A	64[C]	PIA	CB2-CA2	-2.87	1.32	1.35
1	A	64[C]	PIA	CD1-CG2	2.86	1.45	1.39
1	A	64[A]	PIA	CE1-CZ	-2.84	1.33	1.38
1	A	64[C]	PIA	CE1-CZ	-2.60	1.34	1.38
1	A	64[C]	PIA	C1-N3	2.24	1.40	1.37
1	A	64[C]	PIA	CA2-N2	2.22	1.43	1.38
1	A	64[A]	PIA	CA3-N3	-2.22	1.42	1.47

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	64[A]	PIA	O2-C2-CA2	7.47	135.15	130.96
1	A	64[C]	PIA	CB2-CA2-C2	7.40	131.11	122.28
1	A	64[A]	PIA	C2-CA2-N2	-7.08	103.97	108.93
1	A	64[A]	PIA	CA2-N2-C1	6.63	110.66	105.77
1	A	64[A]	PIA	C2-N3-C1	5.94	110.97	107.97
1	A	64[C]	PIA	CB2-CA2-N2	-5.77	120.83	128.83
1	A	64[A]	PIA	CG2-CB2-CA2	-5.19	123.58	129.94
1	A	64[C]	PIA	CD2-CE2-CZ	4.80	125.14	119.88
1	A	64[A]	PIA	CB2-CA2-C2	4.10	127.17	122.28
1	A	64[A]	PIA	CD2-CE2-CZ	4.04	124.31	119.88
1	A	64[A]	PIA	O3-C3-CA3	-3.26	116.56	126.39
1	A	64[A]	PIA	CD2-CG2-CB2	3.18	132.06	121.22
1	A	64[C]	PIA	CA2-N2-C1	3.08	108.04	105.77
1	A	64[C]	PIA	C2-N3-C1	3.07	109.52	107.97
1	A	64[C]	PIA	CD2-CG2-CB2	2.96	131.31	121.22
1	A	64[A]	PIA	CA3-N3-C1	2.83	130.56	127.16
1	A	64[A]	PIA	CE1-CD1-CG2	2.62	124.66	121.25
1	A	64[C]	PIA	OH-CZ-CE1	2.54	127.27	120.02
1	A	64[C]	PIA	CE2-CD2-CG2	-2.45	118.06	121.25
1	A	64[A]	PIA	CA3-N3-C2	-2.42	118.26	123.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	64[C]	PIA	O3-C3-CA3	-2.39	119.18	126.39
1	A	64[C]	PIA	CD1-CG2-CB2	-2.34	113.26	121.22
1	A	64[A]	PIA	CE2-CD2-CG2	-2.32	118.22	121.25
1	A	64[A]	PIA	CD1-CG2-CB2	-2.30	113.39	121.22
1	A	64[A]	PIA	CD2-CG2-CD1	-2.27	114.28	117.64

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	64[A]	PIA	N2-CA2-CB2-CG2
1	A	64[A]	PIA	C2-CA2-CB2-CG2
1	A	64[C]	PIA	C2-CA2-CB2-CG2
1	A	64[A]	PIA	CA2-CB2-CG2-CD1
1	A	64[A]	PIA	CA2-CB2-CG2-CD2
1	A	64[C]	PIA	N2-CA2-CB2-CG2
1	A	64[A]	PIA	N2-C1-CA1-CB1
1	A	64[C]	PIA	C3-CA3-N3-C2
1	A	64[A]	PIA	N3-C1-CA1-CB1

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	64[C]	PIA	2	0
1	A	64[A]	PIA	2	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand 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	$\# Z  > 2$
2	SO4	A	301	-	4,4,4	0.18	0	6,6,6	0.65	0

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 [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, 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	219/220 (99%)	0.22	7 (3%)	47 39	5, 8, 22, 79	1 (0%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	222	ASN	9.8
1	A	1	MET	8.6
1	A	182	LYS	3.9
1	A	2	SER	3.7
1	A	221	PRO	3.3
1	A	150	ASP	3.3
1	A	100[A]	ILE	3.0

### 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, 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
1	PIA	A	64[A]	20/21	0.99	0.10	1,8,12,14	33
1	PIA	A	64[C]	20/21	0.99	0.10	1,5,13,21	33

### 6.3 Carbohydrates

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( $\text{\AA}^2$ )	Q<0.9
2	SO4	A	301	5/5	0.91	0.19	27,28,30,38	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.