Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 10:58 PM EDT

PDB ID	:	5WWN
Title	:	Crystal structure of Tsr1
Authors	:	Ye, K.; Wang, B.
Deposited on	:	2017-01-03
Resolution	:	2.81 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 (i) 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 (i)) were used in the production of this report:

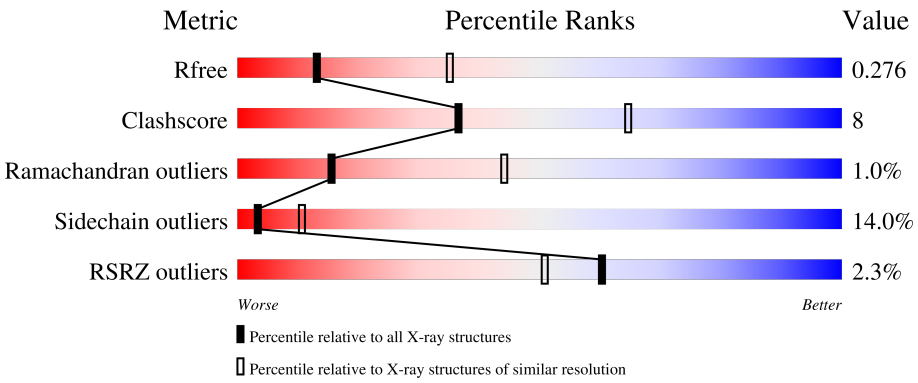
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 <=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	709	2% 53% 18% • 25%

2 Entry composition (i)

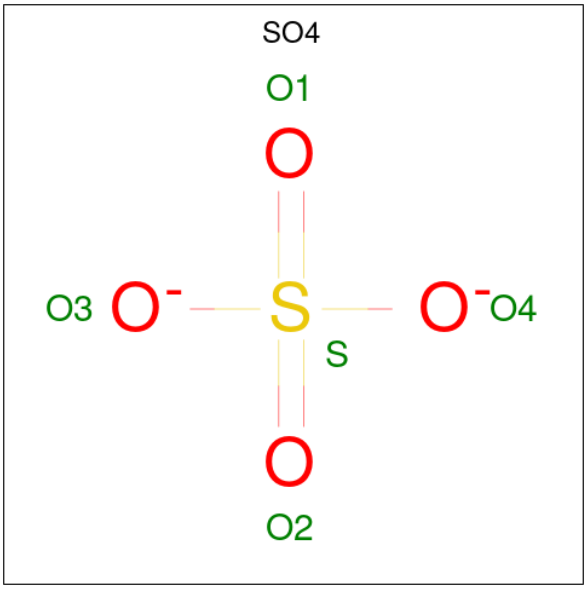
There are 2 unique types of molecules in this entry. The entry contains 4397 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ribosome biogenesis protein TSR1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	531	4347	2801	752	781	13	0	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

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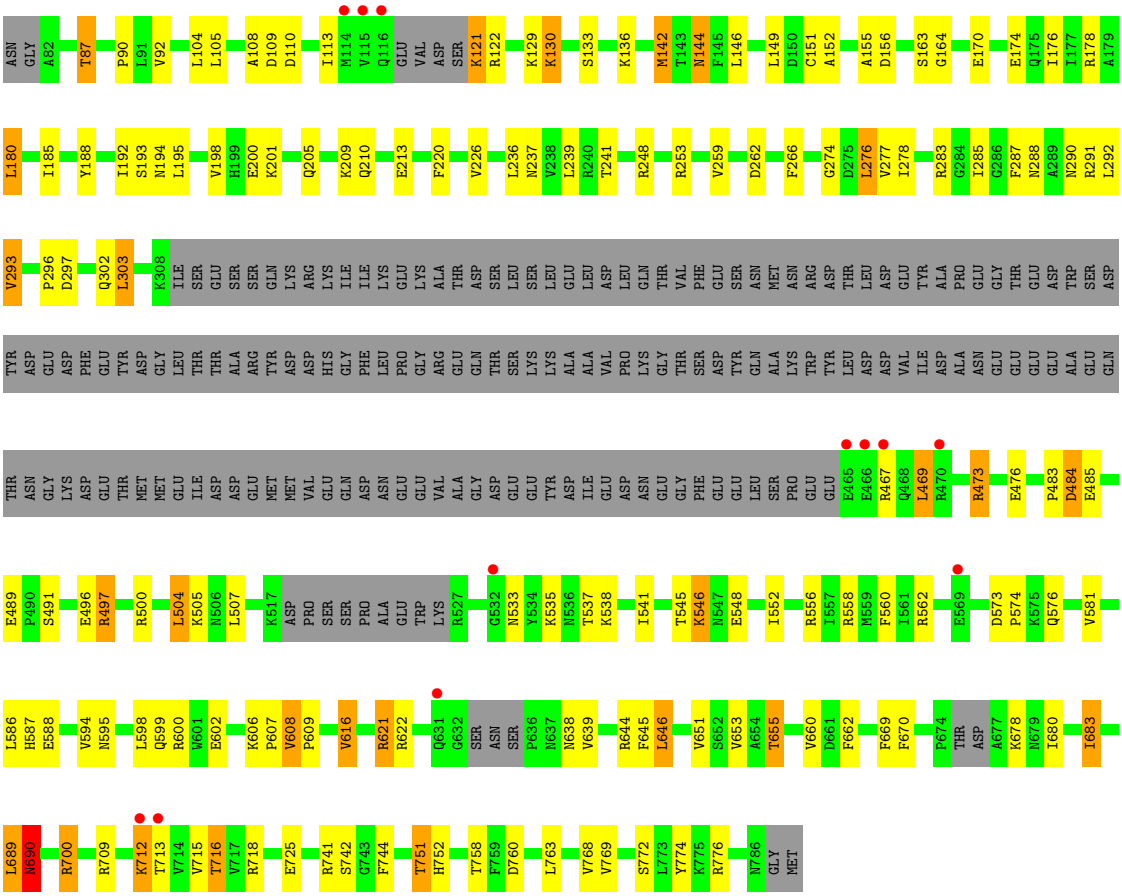
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O	S		
2	A	1	5	4	1	0	0

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: Ribosome biogenesis protein TSR1

Chain A: 2% 53% 18% • 25%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	65.37Å 116.34Å 119.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.46 – 2.81 43.46 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.4 (43.46-2.81) 99.4 (43.46-2.81)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.35 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.214 , 0.274 0.217 , 0.276	Depositor DCC
R_{free} test set	1129 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å^2)	54.6	Xtriage
Anisotropy	0.123	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 46.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.009 for -h,l,k	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	4397	wwPDB-VP
Average B, all atoms (Å^2)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.81% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: 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	0.48	0/4446	0.63	0/5999

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	193	SER	Peptide
1	A	689	LEU	Peptide
1	A	690	ASN	Peptide

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	4347	0	4355	69	0
2	A	50	0	0	1	0
All	All	4397	0	4355	69	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 8.

All (69) 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:152:ALA:HB1	1:A:180:LEU:HD11	1.61	0.81
1:A:497:ARG:NH2	2:A:801:SO4:O4	2.15	0.79
1:A:700:ARG:HG3	1:A:774:TYR:CZ	2.21	0.76
1:A:180:LEU:HD12	1:A:185:ILE:HD12	1.71	0.72
1:A:588:GLU:O	1:A:621:ARG:NH2	2.24	0.70
1:A:164:GLY:H	1:A:194:ASN:HD22	1.41	0.69
1:A:600:ARG:NH2	1:A:606:LYS:O	2.25	0.68
1:A:87:THR:HG23	1:A:155:ALA:HB2	1.80	0.63
1:A:297:ASP:OD1	1:A:622:ARG:NH2	2.35	0.60
1:A:609:PRO:HB3	1:A:645:PHE:CE2	2.38	0.59
1:A:152:ALA:CB	1:A:180:LEU:HD11	2.31	0.58
1:A:213:GLU:HG3	1:A:226:VAL:HB	1.85	0.58
1:A:198:VAL:O	1:A:205:GLN:NE2	2.35	0.57
1:A:285:ILE:HD11	1:A:546:LYS:HB2	1.86	0.56
1:A:715:VAL:HG21	1:A:763:LEU:HD12	1.86	0.56
1:A:108:ALA:O	1:A:110:ASP:N	2.39	0.55
1:A:290:ASN:O	1:A:291:ARG:NH2	2.40	0.54
1:A:110:ASP:OD2	1:A:129:LYS:HE2	2.09	0.53
1:A:709:ARG:HB2	1:A:716:THR:HG23	1.90	0.52
1:A:87:THR:HG21	1:A:151:CYS:O	2.11	0.50
1:A:259:VAL:HG23	1:A:287:PHE:CE2	2.47	0.50
1:A:595:ASN:OD1	1:A:655:THR:HB	2.11	0.50
1:A:545:THR:HA	1:A:548:GLU:HG2	1.93	0.49
1:A:277:VAL:HG22	1:A:560:PHE:CD2	2.49	0.48
1:A:288:ASN:HD21	1:A:290:ASN:HB2	1.78	0.47
1:A:751:THR:HG22	1:A:752:HIS:CE1	2.49	0.47
1:A:469:LEU:HD13	1:A:473:ARG:HG3	1.95	0.47
1:A:538:LYS:HG3	1:A:662:PHE:O	2.14	0.47
1:A:476:GLU:OE2	1:A:607:PRO:HB3	2.15	0.46
1:A:598:LEU:HD13	1:A:669:PHE:CD1	2.51	0.46
1:A:653:VAL:HG12	1:A:655:THR:HG22	1.97	0.46
1:A:285:ILE:HD11	1:A:546:LYS:HE3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:533:ASN:O	1:A:537:THR:HG23	2.15	0.46
1:A:599:GLN:HB2	1:A:651:VAL:HG22	1.97	0.45
1:A:266:PHE:CZ	1:A:274:GLY:HA3	2.51	0.45
1:A:712:LYS:HE3	1:A:712:LYS:HB2	1.81	0.45
1:A:288:ASN:ND2	1:A:290:ASN:HB2	2.32	0.45
1:A:176:ILE:O	1:A:180:LEU:HB2	2.15	0.44
1:A:188:TYR:CE1	1:A:220:PHE:HB3	2.51	0.44
1:A:136:LYS:NZ	1:A:262:ASP:OD1	2.51	0.44
1:A:507:LEU:HG	1:A:639:VAL:HG23	1.99	0.44
1:A:174:GLU:OE1	1:A:178:ARG:NH1	2.51	0.44
1:A:276:LEU:HD23	1:A:276:LEU:HA	1.82	0.44
1:A:92:VAL:HG21	1:A:163:SER:HB3	1.99	0.44
1:A:616:VAL:O	1:A:622:ARG:HA	2.18	0.43
1:A:504:LEU:HD23	1:A:504:LEU:HA	1.85	0.43
1:A:121:LYS:HD2	1:A:576:GLN:HE21	1.83	0.43
1:A:144:ASN:C	1:A:144:ASN:HD22	2.22	0.43
1:A:483:PRO:HA	1:A:484:ASP:HA	1.76	0.42
1:A:283:ARG:HH11	1:A:283:ARG:HB3	1.84	0.42
1:A:156:ASP:OD1	1:A:283:ARG:NH2	2.52	0.42
1:A:292:LEU:CD2	1:A:302:GLN:HG2	2.49	0.42
1:A:670:PHE:HA	1:A:683:ILE:HG23	2.01	0.42
1:A:293:VAL:HG13	1:A:581:VAL:HG13	2.02	0.41
1:A:303:LEU:HA	1:A:303:LEU:HD12	1.85	0.41
1:A:195:LEU:HD23	1:A:195:LEU:HA	1.81	0.41
1:A:537:THR:O	1:A:541:ILE:HG13	2.20	0.41
1:A:586:LEU:O	1:A:587:HIS:HB2	2.20	0.41
1:A:146:LEU:HD23	1:A:146:LEU:HA	1.86	0.41
1:A:105:LEU:HD22	1:A:113:ILE:HG21	2.03	0.41
1:A:236:LEU:HD23	1:A:236:LEU:HA	1.92	0.41
1:A:296:PRO:HA	1:A:297:ASP:HA	1.78	0.41
1:A:541:ILE:O	1:A:545:THR:HG23	2.21	0.41
1:A:573:ASP:HA	1:A:574:PRO:HD2	1.91	0.41
1:A:500:ARG:NH1	1:A:776:ARG:O	2.35	0.41
1:A:689:LEU:HB2	1:A:690:ASN:HB2	2.03	0.41
1:A:484:ASP:HB3	1:A:644:ARG:NH1	2.36	0.40
1:A:90:PRO:HD2	1:A:142:MET:HE1	2.03	0.40
1:A:608:VAL:HG13	1:A:646:LEU:HB3	2.04	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	519/709 (73%)	473 (91%)	41 (8%)	5 (1%)	15 44

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	109	ASP
1	A	130	LYS
1	A	760	ASP
1	A	712	LYS
1	A	638	ASN

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	477/636 (75%)	410 (86%)	67 (14%)	3 11

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	87	THR
1	A	104	LEU
1	A	121	LYS
1	A	122	ARG
1	A	130	LYS
1	A	133	SER

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Mol	Chain	Res	Type
1	A	142	MET
1	A	144	ASN
1	A	149	LEU
1	A	170	GLU
1	A	180	LEU
1	A	192	ILE
1	A	200	GLU
1	A	201	LYS
1	A	209	LYS
1	A	210	GLN
1	A	237	ASN
1	A	239	LEU
1	A	241	THR
1	A	248	ARG
1	A	253	ARG
1	A	276	LEU
1	A	278	ILE
1	A	293	VAL
1	A	303	LEU
1	A	467	ARG
1	A	469	LEU
1	A	473	ARG
1	A	484	ASP
1	A	485	GLU
1	A	489	GLU
1	A	491	SER
1	A	496	GLU
1	A	497	ARG
1	A	504	LEU
1	A	505	LYS
1	A	535	LYS
1	A	546	LYS
1	A	552	ILE
1	A	556	ARG
1	A	558	ARG
1	A	562	ARG
1	A	594	VAL
1	A	602	GLU
1	A	608	VAL
1	A	616	VAL
1	A	621	ARG
1	A	646	LEU

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Mol	Chain	Res	Type
1	A	655	THR
1	A	660	VAL
1	A	678	LYS
1	A	680	ILE
1	A	683	ILE
1	A	690	ASN
1	A	700	ARG
1	A	713	THR
1	A	716	THR
1	A	718	ARG
1	A	725	GLU
1	A	741	ARG
1	A	742	SER
1	A	744	PHE
1	A	751	THR
1	A	758	THR
1	A	768	VAL
1	A	769	VAL
1	A	772	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	134	ASN
1	A	144	ASN
1	A	194	ASN
1	A	509	ASN
1	A	576	GLN
1	A	591	ASN
1	A	640	HIS
1	A	647	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)

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)

10 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	SO4	A	807	-	4,4,4	0.17	0	6,6,6	0.36	0
2	SO4	A	804	-	4,4,4	0.20	0	6,6,6	0.27	0
2	SO4	A	805	-	4,4,4	0.18	0	6,6,6	0.29	0
2	SO4	A	801	-	4,4,4	0.25	0	6,6,6	0.47	0
2	SO4	A	803	-	4,4,4	0.16	0	6,6,6	0.32	0
2	SO4	A	809	-	4,4,4	0.18	0	6,6,6	0.20	0
2	SO4	A	802	-	4,4,4	0.12	0	6,6,6	0.25	0
2	SO4	A	808	-	4,4,4	0.18	0	6,6,6	0.11	0
2	SO4	A	806	-	4,4,4	0.17	0	6,6,6	0.07	0
2	SO4	A	810	-	4,4,4	0.16	0	6,6,6	0.11	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	SO4	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	531/709 (74%)	0.05	12 (2%) 60 51	29, 44, 77, 111	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	465	GLU	5.6
1	A	466	GLU	3.3
1	A	116	GLN	3.3
1	A	467	ARG	3.1
1	A	470	ARG	3.0
1	A	712	LYS	2.9
1	A	114	MET	2.8
1	A	713	THR	2.3
1	A	115	VAL	2.3
1	A	532	GLY	2.2
1	A	631	GLN	2.2
1	A	569	GLU	2.1

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, 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
2	SO4	A	809	5/5	0.81	0.26	76,77,101,105	0
2	SO4	A	810	5/5	0.81	0.66	205,206,209,210	0
2	SO4	A	807	5/5	0.86	0.52	76,77,87,89	0
2	SO4	A	806	5/5	0.88	0.47	100,100,102,104	0
2	SO4	A	805	5/5	0.89	0.37	79,81,85,91	0
2	SO4	A	803	5/5	0.90	0.15	78,78,84,85	0
2	SO4	A	808	5/5	0.91	0.23	77,81,97,98	0
2	SO4	A	801	5/5	0.92	0.16	52,54,65,67	0
2	SO4	A	804	5/5	0.93	0.46	76,76,86,89	0
2	SO4	A	802	5/5	0.95	0.18	62,64,66,74	0

6.5 Other polymers (i)

There are no such residues in this entry.