



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:30 PM UTC

PDB ID : 4RLV / pdb_00004rlv
Title : Crystal Structure of AnkB 24 Ankyrin Repeats in Complex with AnkR Autoinhibition Segment
Authors : Wei, Z.; Wang, C.; Zhang, M.
Deposited on : 2014-10-18
Resolution : 3.49 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

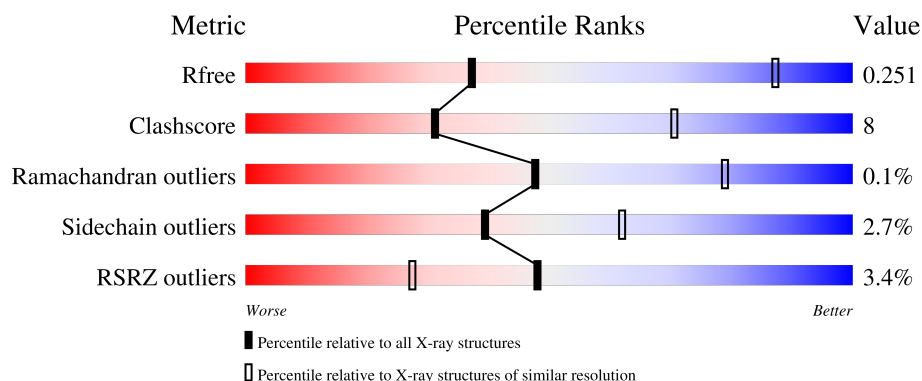
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

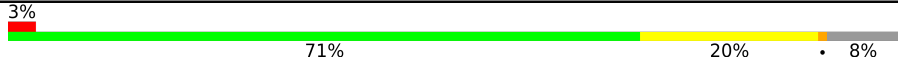
The reported resolution of this entry is 3.49 Å.

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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	910	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6305 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 Ankyrin-1, Ankyrin-2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	838	Total	C	N	O	S	Se	0	0	0
			6260	3905	1156	1176	7	16			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1571	GLY	-	expression tag	UNP D3YTV8
A	1572	PRO	-	expression tag	UNP D3YTV8
A	1573	GLY	-	expression tag	UNP D3YTV8
A	1574	SER	-	expression tag	UNP D3YTV8
A	1575	GLU	-	expression tag	UNP D3YTV8
A	1576	PHE	-	expression tag	UNP D3YTV8
A	1625	GLY	-	linker	UNP D3YTV8
A	1626	SER	-	linker	UNP D3YTV8
A	1627	LEU	-	linker	UNP D3YTV8
A	1628	VAL	-	linker	UNP D3YTV8
A	1629	PRO	-	linker	UNP D3YTV8
A	1630	ARG	-	linker	UNP D3YTV8
A	2024	GLY	-	linker	UNP Q01484
A	2025	SER	-	linker	UNP Q01484
A	2026	GLY	-	linker	UNP Q01484
A	2027	SER	-	linker	UNP Q01484

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	179.69Å 179.69Å 304.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.59 – 3.49 38.59 – 3.49	Depositor EDS
% Data completeness (in resolution range)	94.5 (38.59-3.49) 74.2 (38.59-3.49)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 3.48Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.220 , 0.253 0.221 , 0.251	Depositor DCC
R_{free} test set	1462 reflections (6.28%)	wwPDB-VP
Wilson B-factor (Å ²)	100.4	Xtriage
Anisotropy	0.613	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 87.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6305	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% 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

5.1 Standard geometry

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	0.76	2/6341 (0.0%)	1.09	18/8585 (0.2%)

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	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1591	ILE	CA-CB	5.33	1.61	1.54
1	A	2323	ALA	CA-C	-5.11	1.48	1.53

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1609	VAL	CB-CA-C	-8.99	101.52	111.23
1	A	2189	GLY	N-CA-C	-7.68	106.22	114.67
1	A	2192	ARG	N-CA-C	7.27	119.29	111.36
1	A	1603	GLU	N-CA-C	-6.53	103.74	112.94
1	A	1615	THR	CA-C-N	6.14	126.15	119.89
1	A	1615	THR	C-N-CA	6.14	126.15	119.89
1	A	1618	LEU	N-CA-C	5.89	118.84	109.24
1	A	2146	VAL	CB-CA-C	-5.51	104.91	111.97
1	A	2398	GLY	N-CA-C	-5.46	107.78	115.32
1	A	2522	HIS	CA-C-N	5.46	125.55	119.87
1	A	2522	HIS	C-N-CA	5.46	125.55	119.87

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2477	VAL	N-CA-C	5.46	115.66	110.42
1	A	2490	VAL	N-CA-C	5.25	115.98	110.62
1	A	1596	THR	N-CA-C	-5.18	107.09	113.41
1	A	2437	VAL	N-CA-C	5.11	115.32	110.42
1	A	2323	ALA	CA-C-N	5.11	126.22	119.84
1	A	2323	ALA	C-N-CA	5.11	126.22	119.84
1	A	2357	PRO	CA-C-O	-5.05	115.97	121.27

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1607	MSE	Peptide

5.2 Too-close contacts ⓘ

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	6260	0	6385	104	0
2	A	45	0	0	3	0
All	All	6305	0	6385	104	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 (104) 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:2583:LEU:HD13	1:A:2618:LYS:HG3	1.66	0.78
1:A:2649:LEU:HD13	1:A:2684:LYS:HG3	1.67	0.76
1:A:2048:VAL:HG13	1:A:2086:ARG:HH21	1.57	0.70
1:A:2757:LYS:HG2	1:A:2763:THR:HG22	1.76	0.67
1:A:2721:GLN:O	1:A:2730:THR:OG1	2.07	0.67
1:A:2369:LEU:HD11	1:A:2381:THR:HG23	1.78	0.66
1:A:2046:LYS:HE2	1:A:2050:TYR:HE2	1.62	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2798:LEU:HD21	1:A:2814:LYS:HE2	1.81	0.63
1:A:2035:PHE:HE1	1:A:2047:VAL:HG13	1.64	0.63
1:A:2253:LEU:HA	1:A:2257:ALA:HB3	1.81	0.62
1:A:2637:ALA:HA	1:A:2677:MSE:HE2	1.81	0.62
1:A:2050:TYR:HD1	1:A:2055:ILE:HG13	1.66	0.61
1:A:2800:ILE:HG23	1:A:2803:ARG:HH21	1.66	0.61
1:A:2253:LEU:HD23	1:A:2257:ALA:HB3	1.83	0.60
1:A:2507:LEU:HB3	1:A:2509:LYS:HE2	1.85	0.57
1:A:2241:ALA:O	1:A:2278:ASN:ND2	2.37	0.57
1:A:2791:THR:H	1:A:2795:ASN:H	1.50	0.57
1:A:2601:HIS:HD2	1:A:2635:ILE:HD11	1.69	0.56
1:A:2050:TYR:CD1	1:A:2055:ILE:HG13	2.40	0.56
1:A:2138:ALA:HB1	1:A:2170:ALA:HB2	1.87	0.55
1:A:2481:ARG:NH2	2:A:2907:SO4:O4	2.35	0.55
1:A:2189:GLY:HA2	1:A:2192:ARG:HB2	1.90	0.54
1:A:2427:ILE:HD12	1:A:2431:GLY:HA2	1.88	0.54
1:A:2330:LYS:HG2	1:A:2331:ASN:N	2.23	0.53
1:A:2270:LEU:HD11	1:A:2282:VAL:HG13	1.91	0.53
1:A:2703:ALA:HB2	1:A:2743:MSE:HE2	1.90	0.53
1:A:1607:MSE:HE2	1:A:2267:ILE:HD12	1.91	0.52
1:A:2682:LEU:HD23	1:A:2686:ALA:HB3	1.93	0.51
1:A:2723:ALA:O	1:A:2730:THR:OG1	2.28	0.50
1:A:2567:LEU:HD23	1:A:2599:PRO:HG2	1.93	0.50
1:A:2724:HIS:HB3	1:A:2728:GLY:HA2	1.93	0.50
1:A:2465:GLU:OE2	1:A:2473:ARG:NH2	2.45	0.50
1:A:2061:ASN:HD21	1:A:2065:LEU:HD12	1.76	0.50
1:A:2328:ARG:HB3	1:A:2332:GLY:HA2	1.94	0.50
1:A:2329:THR:HG21	1:A:2338:MSE:CE	2.42	0.49
1:A:2193:LEU:HB3	1:A:2194:PRO:HD3	1.94	0.49
1:A:2640:ASN:HB2	1:A:2674:HIS:CE1	2.48	0.48
1:A:2141:ASN:HB2	1:A:2175:HIS:CD2	2.48	0.48
1:A:2769:ALA:HA	1:A:2777:ILE:HD11	1.94	0.48
1:A:1585:VAL:HG21	1:A:2470:MSE:HE1	1.95	0.48
1:A:2144:ASP:N	1:A:2144:ASP:OD1	2.45	0.48
1:A:2514:GLN:HG2	1:A:2548:VAL:HG11	1.95	0.48
1:A:1596:THR:O	1:A:1600:THR:HG23	2.14	0.48
1:A:2212:LEU:HD23	1:A:2225:MSE:HE2	1.96	0.48
1:A:2329:THR:HG21	1:A:2338:MSE:HE2	1.95	0.48
1:A:2343:ASP:OD2	1:A:2377:HIS:HB3	2.14	0.47
1:A:2508:GLY:HA2	1:A:2545:VAL:HG21	1.96	0.47
1:A:1599:ASP:OD2	1:A:2330:LYS:NZ	2.48	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2232:SER:HB2	1:A:2264:ARG:HH21	1.78	0.47
1:A:1625:GLY:O	1:A:1627:LEU:HG	2.14	0.46
1:A:2766:HIS:NE2	1:A:2795:ASN:O	2.47	0.46
1:A:2691:SER:HB2	1:A:2695:GLY:HA2	1.97	0.46
1:A:2084:LEU:HD13	1:A:2119:GLU:HG2	1.97	0.46
1:A:2787:PRO:HG3	1:A:2816:VAL:O	2.15	0.46
1:A:1597:GLU:O	1:A:1600:THR:OG1	2.32	0.46
1:A:2303:LEU:HD11	1:A:2315:VAL:HG13	1.97	0.46
1:A:2798:LEU:HD11	1:A:2814:LYS:HE3	1.98	0.46
1:A:1588:ILE:HD11	1:A:2437:VAL:HG13	1.96	0.45
1:A:2295:LYS:HE2	1:A:2295:LYS:HB3	1.73	0.45
1:A:2385:LEU:HD13	1:A:2420:TYR:CG	2.51	0.45
1:A:2572:LYS:HG3	1:A:2606:TYR:CE2	2.51	0.45
1:A:2167:LEU:HD23	1:A:2191:VAL:HG12	1.99	0.45
1:A:2079:LEU:O	1:A:2083:LEU:HG	2.17	0.45
1:A:1613:GLY:C	1:A:1614:LEU:HD23	2.42	0.44
1:A:2202:LYS:HE3	2:A:2909:SO4:O3	2.17	0.44
1:A:2411:ILE:HG12	1:A:2446:ILE:HD11	1.99	0.44
1:A:2336:LEU:HD11	1:A:2348:VAL:HG13	1.99	0.44
1:A:2550:LEU:HD13	1:A:2585:ARG:HG3	2.00	0.44
1:A:2316:GLU:HB2	1:A:2350:HIS:CD2	2.52	0.44
1:A:2697:THR:O	1:A:2700:HIS:HB2	2.18	0.43
1:A:2690:MSE:HE2	1:A:2690:MSE:HB3	1.96	0.43
1:A:2262:THR:HG22	1:A:2268:THR:HG22	2.00	0.43
1:A:2770:GLN:HG2	1:A:2771:GLN:NE2	2.33	0.43
1:A:2134:LEU:HD21	1:A:2156:GLN:HB3	2.00	0.43
1:A:2427:ILE:HG22	1:A:2433:THR:HG22	2.00	0.43
1:A:2507:LEU:HB2	1:A:2509:LYS:HG2	2.00	0.43
1:A:1601:MSE:O	1:A:1605:SER:N	2.52	0.43
1:A:2162:ASP:OD2	1:A:2162:ASP:N	2.49	0.43
1:A:2083:LEU:HD23	1:A:2083:LEU:HA	1.74	0.42
1:A:2277:GLY:HA2	1:A:2314:VAL:HG21	2.01	0.42
1:A:2343:ASP:HB2	1:A:2377:HIS:CD2	2.55	0.42
1:A:2495:ARG:O	1:A:2498:GLN:HG2	2.20	0.42
1:A:2676:ASP:N	1:A:2676:ASP:OD1	2.51	0.42
1:A:2732:LEU:HD22	1:A:2752:ALA:HB1	2.01	0.42
1:A:2188:LYS:O	1:A:2192:ARG:HG3	2.20	0.41
1:A:2357:PRO:HB2	1:A:2360:ASP:HB2	2.03	0.41
1:A:2145:VAL:O	1:A:2149:LEU:HG	2.20	0.41
1:A:2754:VAL:HA	1:A:2764:PRO:HG2	2.02	0.41
1:A:2300:LEU:HA	1:A:2300:LEU:HD23	1.76	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1597:GLU:HG3	1:A:1598:HIS:N	2.30	0.41
1:A:2310:GLY:HA2	1:A:2347:CYS:SG	2.60	0.41
1:A:2343:ASP:HA	1:A:2380:VAL:HG21	2.02	0.41
1:A:2129:ASN:HB2	1:A:2131:PHE:CE2	2.55	0.41
1:A:2402:LEU:HD12	1:A:2402:LEU:HA	1.73	0.41
1:A:2692:THR:HG23	1:A:2696:LEU:O	2.21	0.41
1:A:2746:PHE:HA	1:A:2749:LYS:HD2	2.02	0.41
1:A:2227:ASN:HB3	1:A:2261:PHE:CE1	2.56	0.41
1:A:2673:GLY:HA2	1:A:2710:VAL:HG21	2.03	0.41
1:A:2601:HIS:CD2	1:A:2635:ILE:HD11	2.53	0.41
1:A:2688:ILE:H	1:A:2688:ILE:HG13	1.62	0.41
1:A:2244:GLY:HA2	1:A:2281:MSE:HG3	2.03	0.40
1:A:2436:HIS:HD2	1:A:2470:MSE:HG3	1.86	0.40
1:A:1594:ALA:HB2	2:A:2908:SO4:O1	2.21	0.40
1:A:2378:TYR:CE1	1:A:2416:LEU:HD22	2.56	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	834/910 (92%)	791 (95%)	42 (5%)	1 (0%)	48 79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1610	TRP

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	660/706 (94%)	642 (97%)	18 (3%)	39 62

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1585	VAL
1	A	1587	VAL
1	A	1597	GLU
1	A	1601	MSE
1	A	1611	SER
1	A	1614	LEU
1	A	1616	PRO
1	A	1622	GLU
1	A	1628	VAL
1	A	2162	ASP
1	A	2171	LEU
1	A	2172	GLN
1	A	2192	ARG
1	A	2330	LYS
1	A	2470	MSE
1	A	2595	ASN
1	A	2730	THR
1	A	2734	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1608	GLN
1	A	2129	ASN
1	A	2173	GLN
1	A	2595	ASN
1	A	2671	GLN
1	A	2704	GLN
1	A	2740	ASN

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](#)

9 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	2909	-	4,4,4	0.24	0	6,6,6	0.14	0
2	SO4	A	2906	-	4,4,4	0.29	0	6,6,6	0.33	0
2	SO4	A	2904	-	4,4,4	0.25	0	6,6,6	0.27	0
2	SO4	A	2901	-	4,4,4	0.20	0	6,6,6	0.52	0
2	SO4	A	2903	-	4,4,4	0.22	0	6,6,6	0.23	0
2	SO4	A	2908	-	4,4,4	0.29	0	6,6,6	0.23	0
2	SO4	A	2905	-	4,4,4	0.27	0	6,6,6	0.41	0
2	SO4	A	2907	-	4,4,4	0.30	0	6,6,6	0.21	0
2	SO4	A	2902	-	4,4,4	0.35	0	6,6,6	0.19	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.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2909	SO4	1	0
2	A	2908	SO4	1	0
2	A	2907	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

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	822/910 (90%)	0.20	28 (3%)	48 27	28, 108, 225, 265	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2769	ALA	7.3
1	A	2776	ILE	4.9
1	A	1591	ILE	4.1
1	A	1590	ALA	3.9
1	A	2755	ASN	3.7
1	A	2813	LEU	3.6
1	A	2816	VAL	3.2
1	A	2789	ALA	3.1
1	A	2765	LEU	3.1
1	A	2768	ALA	3.0
1	A	2736	CYS	2.8
1	A	2787	PRO	2.6
1	A	2790	THR	2.6
1	A	1587	VAL	2.5
1	A	2815	VAL	2.3
1	A	1585	VAL	2.2
1	A	2732	LEU	2.2
1	A	2764	PRO	2.2
1	A	2748	LEU	2.1
1	A	2794	GLY	2.1
1	A	2699	LEU	2.1
1	A	1593	LEU	2.1
1	A	2770	GLN	2.1
1	A	1596	THR	2.0
1	A	2792	ALA	2.0
1	A	1588	ILE	2.0
1	A	2671	GLN	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	2780	LEU	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(Å ²)	Q<0.9
2	SO4	A	2908	5/5	0.39	0.14	140,143,146,150	0
2	SO4	A	2909	5/5	0.70	0.12	137,138,140,142	0
2	SO4	A	2905	5/5	0.73	0.08	121,125,128,130	0
2	SO4	A	2901	5/5	0.79	0.08	92,112,116,116	0
2	SO4	A	2902	5/5	0.80	0.10	93,100,102,104	0
2	SO4	A	2903	5/5	0.84	0.06	100,101,110,113	0
2	SO4	A	2904	5/5	0.86	0.08	108,112,114,116	5
2	SO4	A	2906	5/5	0.88	0.06	116,117,126,133	0
2	SO4	A	2907	5/5	0.88	0.07	142,143,145,148	0

6.5 Other polymers [i](#)

There are no such residues in this entry.