



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 04:23 PM UTC

PDB ID : 1OLO / pdb_00001olo
Title : Hexameric Replicative DNA Helicase RepA from Plasmid RSF1010 - Cubic
Crystal Structure
Authors : Niedenzu, T.; Saenger, W.
Deposited on : 2003-08-08
Resolution : 2.55 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

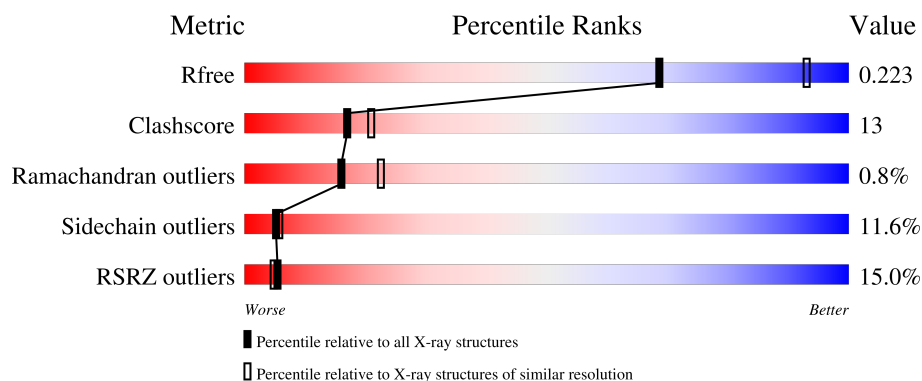
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	279	9% 68% 18% • 11%
1	B	279	18% 71% 14% • 10%

2 Entry composition [i](#)

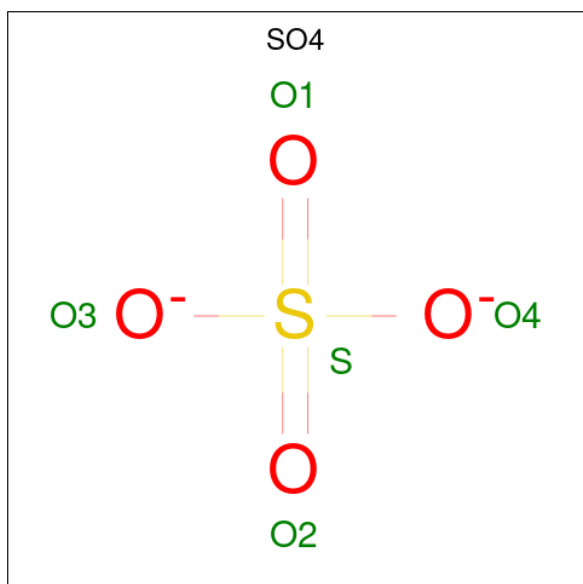
There are 3 unique types of molecules in this entry. The entry contains 4295 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 REGULATORY PROTEIN REPA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	7	0
			1941	1232	352	349	8			
1	B	250	Total	C	N	O	S	0	3	0
			1923	1221	349	345	8			

- 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	B	1	Total	O	S	0	0
			5	4	1		

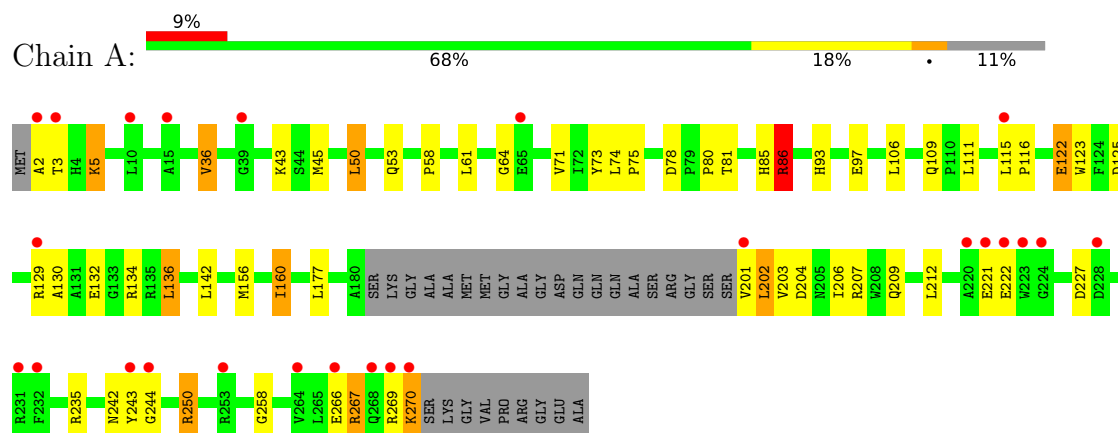
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	212	Total 212	O 212	0	0
3	B	209	Total 209	O 209	0	0

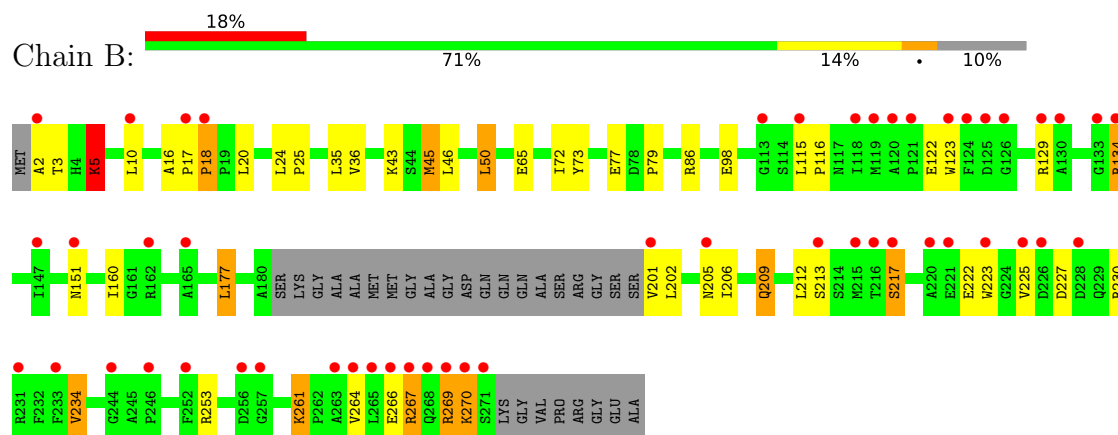
3 Residue-property plots [i](#)

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: REGULATORY PROTEIN REPA



• Molecule 1: REGULATORY PROTEIN REPA



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 3 2	Depositor
Cell constants a, b, c, α , β , γ	190.29Å 190.29Å 190.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.55 25.00 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.8 (25.00-2.55) 99.7 (25.00-2.55)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.176 , 0.211 0.191 , 0.223	Depositor DCC
R_{free} test set	991 reflections (2.48%)	wwPDB-VP
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4295	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% 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: 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	1.10	1/2016 (0.0%)	1.12	4/2738 (0.1%)
1	B	1.15	3/1984 (0.2%)	1.11	2/2695 (0.1%)
All	All	1.12	4/4000 (0.1%)	1.12	6/5433 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	18	PRO	CA-C	7.00	1.58	1.52
1	A	156	MET	SD-CE	-6.76	1.62	1.79
1	B	79	PRO	CA-C	5.23	1.55	1.52
1	B	5	LYS	CG-CD	5.09	1.67	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	234	VAL	CB-CA-C	-7.61	97.55	110.71
1	A	36	VAL	N-CA-C	6.68	118.11	108.48
1	A	86	ARG	N-CA-C	-6.02	104.64	111.14
1	A	64	GLY	CA-C-O	-5.96	118.12	122.23
1	A	204	ASP	N-CA-C	5.27	118.49	111.75
1	B	5	LYS	CG-CD-CE	5.04	122.90	111.30

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	1941	0	1923	49	1
1	B	1923	0	1922	50	2
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	212	0	0	19	0
3	B	209	0	0	19	2
All	All	4295	0	3845	98	3

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 13.

All (98) 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:244:GLY:HA3	3:A:2025:HOH:O	1.35	1.21
1:A:45:MET:HE1	1:A:86:ARG:HG3	1.41	0.98
1:A:45:MET:CE	1:A:86:ARG:HG3	1.95	0.96
1:B:73:TYR:CE1	3:B:2097:HOH:O	2.18	0.94
1:B:201:VAL:HG21	3:B:2107:HOH:O	1.71	0.90
1:B:201:VAL:HG23	1:B:202:LEU:N	1.90	0.86
1:A:266:GLU:HB3	3:A:2207:HOH:O	1.75	0.86
1:B:43:LYS:HB3	1:B:177:LEU:HD13	1.61	0.83
1:A:43:LYS:HB3	1:A:177:LEU:HD13	1.61	0.82
1:B:201:VAL:HG23	1:B:202:LEU:H	1.44	0.82
1:A:53[A]:GLN:HG2	1:A:58:PRO:O	1.81	0.81
1:B:261:LYS:HE3	3:B:2206:HOH:O	1.82	0.80
1:B:201:VAL:CG2	3:B:2107:HOH:O	2.27	0.79
1:A:85[B]:HIS:CD2	3:A:2088:HOH:O	2.35	0.78
1:B:3:THR:HB	3:B:2003:HOH:O	1.83	0.77
1:B:5:LYS:N	1:B:5:LYS:HD2	2.01	0.76
1:A:244:GLY:CA	3:A:2025:HOH:O	2.10	0.74
1:B:72:ILE:HB	1:B:134[B]:ARG:HD3	1.68	0.74
1:B:5:LYS:H	1:B:5:LYS:CD	2.01	0.74
1:A:160:ILE:HD11	1:A:202:LEU:HD12	1.67	0.74
1:B:16:ALA:O	3:B:2032:HOH:O	2.05	0.73
1:B:206:ILE:HD11	1:B:209:GLN:HG3	1.71	0.71
1:A:85[B]:HIS:NE2	3:A:2088:HOH:O	2.25	0.70
1:A:221:GLU:OE2	3:A:2173:HOH:O	2.09	0.69
1:B:115:LEU:HD23	3:B:2121:HOH:O	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:253[B]:ARG:NH2	1:B:264:VAL:H	1.90	0.69
1:A:45:MET:CE	1:A:86:ARG:CG	2.73	0.66
1:A:73:TYR:HD2	1:A:75:PRO:HD3	1.61	0.66
1:A:5:LYS:HE3	3:B:2060:HOH:O	1.97	0.65
1:B:5:LYS:HD2	1:B:5:LYS:H	1.57	0.65
1:B:223:TRP:CE3	1:B:267:ARG:HB2	2.33	0.64
1:A:50:LEU:HD13	1:A:61:LEU:CD1	2.28	0.63
1:A:73:TYR:CD2	1:A:75:PRO:HD3	2.34	0.62
1:A:242[A]:ASN:ND2	3:A:2187:HOH:O	2.32	0.62
1:A:160:ILE:HD11	1:A:202:LEU:CD1	2.29	0.62
1:B:2:ALA:HB1	3:B:2002:HOH:O	2.00	0.62
1:B:223:TRP:CZ3	1:B:267:ARG:HB2	2.35	0.62
1:B:213:SER:HB2	3:B:2181:HOH:O	1.99	0.61
1:A:45:MET:HE2	1:A:86:ARG:HG3	1.81	0.61
1:B:201:VAL:HG13	3:B:2175:HOH:O	2.01	0.60
1:A:50:LEU:HD13	1:A:61:LEU:HD12	1.83	0.60
1:B:16:ALA:HB1	1:B:17:PRO:HD2	1.83	0.59
1:A:74:LEU:HD23	1:A:109:GLN:HB3	1.86	0.57
1:A:129[A]:ARG:HD2	3:A:2061:HOH:O	2.03	0.57
1:A:270:LYS:CE	3:A:2175:HOH:O	2.51	0.57
1:B:45:MET:CE	1:B:86:ARG:HB3	2.35	0.56
1:A:129[A]:ARG:CD	3:A:2061:HOH:O	2.54	0.55
1:A:71:VAL:HG13	1:A:136:LEU:HB3	1.89	0.55
1:B:201:VAL:CG2	1:B:202:LEU:N	2.61	0.55
1:A:45:MET:HE1	1:A:86:ARG:CG	2.26	0.54
1:A:242[A]:ASN:CG	3:A:2187:HOH:O	2.51	0.54
1:B:266:GLU:O	1:B:267:ARG:HB3	2.08	0.53
1:A:45:MET:HE2	1:A:86:ARG:CG	2.38	0.52
1:A:93:HIS:HE1	3:A:2200:HOH:O	1.92	0.51
1:A:142:LEU:HD23	1:A:202:LEU:HD21	1.92	0.50
1:A:160:ILE:HG12	1:A:206:ILE:HD11	1.94	0.50
1:B:253[B]:ARG:HH21	1:B:264:VAL:H	1.59	0.50
1:B:160:ILE:HD13	1:B:205:ASN:HB3	1.93	0.49
1:B:35:LEU:HB3	1:B:177:LEU:HD22	1.94	0.49
1:B:72:ILE:CG2	1:B:134[B]:ARG:HD3	2.42	0.49
1:B:45:MET:HE3	1:B:86:ARG:HB3	1.95	0.48
1:B:3:THR:HG22	3:B:2004:HOH:O	2.13	0.48
1:A:116:PRO:HA	1:A:123:TRP:CE2	2.48	0.48
1:B:116:PRO:HA	1:B:123:TRP:CE2	2.48	0.48
1:B:264:VAL:HA	3:B:2207:HOH:O	2.13	0.48
1:B:227:ASP:O	1:B:230:ARG:HG2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:SER:HB2	3:B:2183:HOH:O	2.14	0.47
1:A:5:LYS:HD3	3:A:2004:HOH:O	2.14	0.47
1:A:235:ARG:NH1	3:A:2182:HOH:O	2.47	0.47
1:A:250:ARG:NH2	3:A:2193:HOH:O	2.48	0.46
1:B:98:GLU:HG2	3:B:2082:HOH:O	2.15	0.46
1:A:267:ARG:HA	1:A:267:ARG:HD2	1.56	0.45
1:A:207:ARG:HE	1:A:242[B]:ASN:ND2	2.14	0.45
1:B:72:ILE:CB	1:B:134[B]:ARG:HD3	2.41	0.45
1:A:53[A]:GLN:OE1	3:A:2054:HOH:O	2.21	0.45
1:A:129[A]:ARG:HG2	3:A:2061:HOH:O	2.18	0.44
1:B:24:LEU:HB3	1:B:25:PRO:HD2	1.99	0.44
1:A:130:ALA:O	1:A:134[A]:ARG:HD2	2.18	0.44
1:A:207:ARG:HE	1:A:242[B]:ASN:HD21	1.65	0.44
1:A:207:ARG:HD2	1:B:77:GLU:HB3	2.00	0.44
1:B:5:LYS:HE2	3:B:2008:HOH:O	2.17	0.44
1:B:5:LYS:NZ	3:B:2009:HOH:O	2.45	0.44
1:A:86:ARG:HG2	1:A:86:ARG:HH11	1.83	0.43
1:B:151:ASN:OD1	1:B:151:ASN:C	2.60	0.43
1:A:74:LEU:CD2	1:A:109:GLN:HB3	2.46	0.43
1:A:2:ALA:HB1	3:A:2001:HOH:O	2.17	0.43
1:B:73:TYR:CZ	3:B:2097:HOH:O	2.49	0.43
1:B:201:VAL:CG2	1:B:202:LEU:H	2.14	0.43
1:B:223:TRP:HA	1:B:267:ARG:HA	2.01	0.43
1:A:266:GLU:CB	3:A:2207:HOH:O	2.50	0.42
1:B:206:ILE:CD1	1:B:209:GLN:HG3	2.43	0.42
1:B:177:LEU:N	1:B:177:LEU:HD23	2.34	0.42
1:A:125:ASP:O	1:A:129[B]:ARG:HG2	2.19	0.42
1:A:45:MET:HG2	1:A:258:GLY:O	2.20	0.41
1:B:36:VAL:HG21	1:B:209:GLN:OE1	2.20	0.41
1:B:122:GLU:HB3	3:B:2128:HOH:O	2.21	0.41
1:B:46:LEU:HG	1:B:50:LEU:HD22	2.02	0.41

All (3) 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:85[A]:HIS:NE2	1:B:18:PRO:O[9_555]	1.92	0.28
1:B:18:PRO:CB	3:B:2018:HOH:O[13_545]	2.06	0.14
3:B:2018:HOH:O	3:B:2018:HOH:O[13_545]	2.13	0.07

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	252/279 (90%)	249 (99%)	2 (1%)	1 (0%)	30	39
1	B	249/279 (89%)	241 (97%)	5 (2%)	3 (1%)	10	13
All	All	501/558 (90%)	490 (98%)	7 (1%)	4 (1%)	16	22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	222	GLU
1	B	269	ARG
1	A	243	TYR
1	B	270	LYS

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	200/212 (94%)	171 (86%)	29 (14%)	3	2
1	B	197/212 (93%)	178 (90%)	19 (10%)	8	9
All	All	397/424 (94%)	349 (88%)	48 (12%)	5	5

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	5	LYS

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Mol	Chain	Res	Type
1	A	36	VAL
1	A	50	LEU
1	A	78	ASP
1	A	80	PRO
1	A	81	THR
1	A	86	ARG
1	A	97[A]	GLU
1	A	97[B]	GLU
1	A	106	LEU
1	A	111	LEU
1	A	115	LEU
1	A	122[A]	GLU
1	A	122[B]	GLU
1	A	132	GLU
1	A	136	LEU
1	A	160	ILE
1	A	201	VAL
1	A	202	LEU
1	A	203	VAL
1	A	209	GLN
1	A	212	LEU
1	A	222	GLU
1	A	227	ASP
1	A	250	ARG
1	A	267	ARG
1	A	269	ARG
1	A	270	LYS
1	B	5	LYS
1	B	10	LEU
1	B	20	LEU
1	B	45	MET
1	B	50	LEU
1	B	65	GLU
1	B	129	ARG
1	B	134[A]	ARG
1	B	134[B]	ARG
1	B	177	LEU
1	B	209	GLN
1	B	212	LEU
1	B	217	SER
1	B	225	VAL
1	B	234	VAL

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Mol	Chain	Res	Type
1	B	261	LYS
1	B	267	ARG
1	B	269	ARG
1	B	270	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	HIS
1	A	158	GLN
1	B	93	HIS
1	B	158	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](#)

2 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	300	-	4,4,4	0.43	0	6,6,6	0.74	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	300	-	4,4,4	0.38	0	6,6,6	0.72	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, 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	249/279 (89%)	0.68	25 (10%)	12 12	18, 31, 46, 51	7 (2%)
1	B	250/279 (89%)	1.01	50 (20%)	3 2	14, 32, 43, 54	3 (1%)
All	All	499/558 (89%)	0.84	75 (15%)	5 5	14, 31, 44, 54	10 (2%)

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	271	SER	7.6
1	A	269	ARG	6.0
1	A	243	TYR	5.9
1	B	221	GLU	4.0
1	B	269	ARG	4.0
1	B	266	GLU	3.9
1	B	223	TRP	3.8
1	B	267	ARG	3.7
1	B	215	MET	3.7
1	A	244	GLY	3.7
1	A	268	GLN	3.7
1	A	270	LYS	3.7
1	B	2	ALA	3.5
1	B	268	GLN	3.3
1	A	266	GLU	3.3
1	B	264	VAL	3.2
1	B	17	PRO	3.1
1	B	231	ARG	3.0
1	B	120	ALA	2.9
1	A	264	VAL	2.9
1	B	113	GLY	2.9
1	B	126	GLY	2.8
1	B	115	LEU	2.7
1	A	222	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	2	ALA	2.7
1	A	15	ALA	2.7
1	B	270	LYS	2.6
1	B	216	THR	2.6
1	A	223	TRP	2.6
1	B	147	ILE	2.6
1	A	232	PHE	2.6
1	B	257	GLY	2.5
1	A	220	ALA	2.5
1	B	265	LEU	2.5
1	A	129[A]	ARG	2.5
1	A	224	GLY	2.5
1	B	244	GLY	2.5
1	B	162	ARG	2.5
1	B	124	PHE	2.4
1	A	39	GLY	2.4
1	A	228	ASP	2.4
1	A	231	ARG	2.4
1	A	65	GLU	2.4
1	B	246	PRO	2.4
1	B	256	ASP	2.3
1	B	121	PRO	2.3
1	B	252	PHE	2.3
1	B	201	VAL	2.3
1	B	220	ALA	2.3
1	B	119	MET	2.3
1	A	201	VAL	2.3
1	B	213	SER	2.2
1	A	3	THR	2.2
1	B	134[A]	ARG	2.2
1	B	118	ILE	2.2
1	B	123	TRP	2.2
1	B	233	PHE	2.2
1	B	151	ASN	2.2
1	B	228	ASP	2.2
1	B	205	ASN	2.1
1	B	129	ARG	2.1
1	B	217	SER	2.1
1	B	225	VAL	2.1
1	A	10	LEU	2.1
1	A	115	LEU	2.1
1	B	165	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	133	GLY	2.1
1	A	221	GLU	2.1
1	B	125	ASP	2.1
1	B	226	ASP	2.1
1	B	263	ALA	2.0
1	A	253	ARG	2.0
1	B	10	LEU	2.0
1	B	130	ALA	2.0
1	B	18	PRO	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	SO4	A	300	5/5	0.96	0.08	25,27,32,36	0
2	SO4	B	300	5/5	0.98	0.06	30,30,32,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.