



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 22, 2021 – 10:07 AM JST

PDB ID : 7DVV
Title : Heme sensor protein PefR from *Streptococcus agalactiae* bound to operator DNA (28-mer)
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Deposited on : 2021-01-15
Resolution : 2.49 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.23.1
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)	:	2.23.1

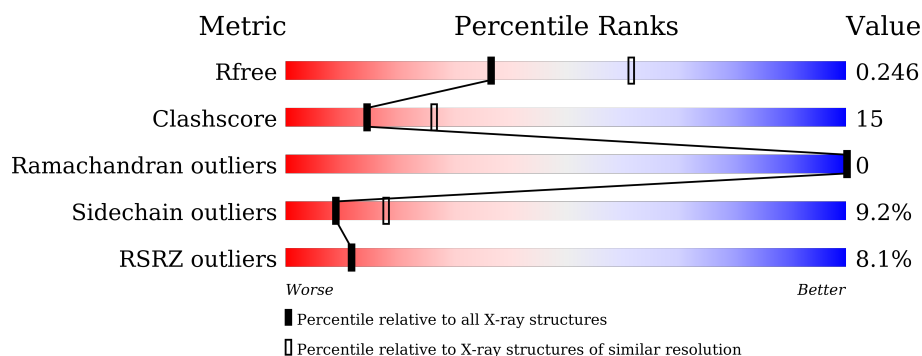
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	153	<div> <div>8%</div> <div>68%</div> <div>16%</div> <div>6%</div> <div>9%</div> </div>
1	B	153	<div> <div>10%</div> <div>63%</div> <div>25%</div> <div>8%</div> </div>
2	L	28	<div> <div>36%</div> <div>57%</div> <div>7%</div> </div>
3	M	28	<div> <div>25%</div> <div>64%</div> <div>11%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3424 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 HTH marR-type domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	139	Total	C	N	O	S	0	0	0
			1129	723	200	201	5			
1	B	140	Total	C	N	O	S	0	0	0
			1138	728	201	204	5			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	147	SER	-	expression tag	UNP Q8E4J9
A	148	LEU	-	expression tag	UNP Q8E4J9
A	149	GLU	-	expression tag	UNP Q8E4J9
A	150	VAL	-	expression tag	UNP Q8E4J9
A	151	LEU	-	expression tag	UNP Q8E4J9
A	152	PHE	-	expression tag	UNP Q8E4J9
A	153	GLN	-	expression tag	UNP Q8E4J9
B	147	SER	-	expression tag	UNP Q8E4J9
B	148	LEU	-	expression tag	UNP Q8E4J9
B	149	GLU	-	expression tag	UNP Q8E4J9
B	150	VAL	-	expression tag	UNP Q8E4J9
B	151	LEU	-	expression tag	UNP Q8E4J9
B	152	PHE	-	expression tag	UNP Q8E4J9
B	153	GLN	-	expression tag	UNP Q8E4J9

- Molecule 2 is a DNA chain called DNA (28-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	28	Total	C	N	O	P	0	0	0
			573	276	105	164	28			

- Molecule 3 is a DNA chain called DNA (28-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	M	28	Total 575	C 278	N 97	O 172	P 28	0	0	0

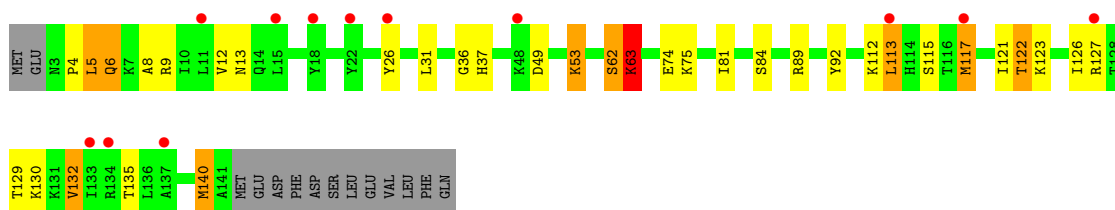
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	O 1	0	0
4	B	1	Total 1	O 1	0	0
4	L	5	Total 5	O 5	0	0
4	M	2	Total 2	O 2	0	0

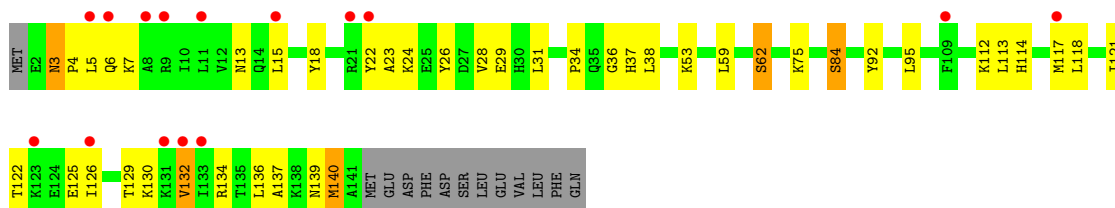
3 Residue-property plots [i](#)

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: HTH marR-type domain-containing protein



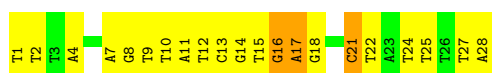
- Molecule 1: HTH marR-type domain-containing protein



- Molecule 2: DNA (28-MER)



- Molecule 3: DNA (28-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 1 2	Depositor
Cell constants a, b, c, α , β , γ	100.97Å 100.97Å 128.99Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.06 – 2.49 47.01 – 2.49	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.06-2.49) 99.9 (47.01-2.49)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.217 , 0.246 0.217 , 0.246	Depositor DCC
R_{free} test set	1262 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	69.0	Xtriage
Anisotropy	0.453	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 69.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.488 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3424	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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](#)

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	0.46	0/1148	0.92	1/1539 (0.1%)
1	B	0.49	0/1157	0.92	0/1551
2	L	0.86	1/643 (0.2%)	1.24	7/989 (0.7%)
3	M	0.96	3/643 (0.5%)	1.22	8/991 (0.8%)
All	All	0.66	4/3591 (0.1%)	1.05	16/5070 (0.3%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	M	24	DT	O3'-P	-9.58	1.49	1.61
2	L	5	DA	O3'-P	-6.12	1.53	1.61
3	M	15	DT	O3'-P	-5.31	1.54	1.61
3	M	16	DG	O3'-P	-5.04	1.55	1.61

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	9	DT	O5'-P-OP1	-12.26	94.66	105.70
3	M	10	DT	O5'-P-OP1	-7.68	98.78	105.70
2	L	17	DA	O5'-P-OP1	7.42	119.61	110.70
3	M	9	DT	O5'-P-OP1	-7.05	99.35	105.70
3	M	21	DC	C1'-O4'-C4'	-6.92	103.18	110.10
3	M	13	DC	C1'-O4'-C4'	-6.64	103.45	110.10
2	L	13	DC	C1'-O4'-C4'	-6.37	103.73	110.10
3	M	4	DA	O5'-P-OP1	-6.24	100.08	105.70
2	L	14	DA	O5'-P-OP1	-5.99	100.31	105.70
3	M	17	DA	O5'-P-OP1	5.87	117.74	110.70
2	L	21	DC	C1'-O4'-C4'	-5.59	104.51	110.10
1	A	63	LYS	CB-CA-C	5.51	121.41	110.40
2	L	17	DA	O5'-P-OP2	-5.46	100.79	105.70
2	L	24	DT	C1'-O4'-C4'	-5.41	104.69	110.10
3	M	14	DG	O5'-P-OP1	-5.11	101.10	105.70
3	M	21	DC	O4'-C4'-C3'	-5.07	102.47	104.50

There are no chirality outliers.

There are no planarity outliers.

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	1129	0	1190	38	0
1	B	1138	0	1196	45	0
2	L	573	0	318	18	0
3	M	575	0	322	17	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	L	5	0	0	0	0
4	M	2	0	0	0	0
All	All	3424	0	3026	96	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 15.

All (96) 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:132:VAL:HG11	1:B:129:THR:HG21	1.38	1.03
1:B:117:MET:HG3	1:B:118:LEU:HD12	1.53	0.88
2:L:17:DA:C2'	2:L:18:DT:H5''	2.05	0.87
2:L:17:DA:H2'	2:L:18:DT:H5''	1.59	0.84
3:M:1:DT:H2''	3:M:2:DT:H71	1.63	0.79
1:B:3:ASN:O	1:B:7:LYS:HG2	1.84	0.77
1:B:84:SER:HB2	1:B:92:TYR:CE2	2.20	0.76
3:M:27:DT:H4'	3:M:28:DA:OP1	1.86	0.75
3:M:1:DT:H2''	3:M:2:DT:C7	2.19	0.72
1:A:122:THR:O	1:A:126:ILE:HG13	1.93	0.69
1:A:121:ILE:HG22	1:A:126:ILE:HG13	1.76	0.68
1:B:125:GLU:O	1:B:129:THR:HG23	1.94	0.68
1:B:117:MET:HG3	1:B:118:LEU:CD1	2.23	0.67
2:L:17:DA:H2''	2:L:18:DT:H5''	1.78	0.66
3:M:7:DA:H2''	3:M:8:DG:H5'	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:MET:HG3	1:B:18:TYR:CD2	2.35	0.62
1:A:113:LEU:HD22	1:B:5:LEU:HD11	1.81	0.62
2:L:11:DC:H2''	2:L:12:DT:H5'	1.82	0.62
1:A:121:ILE:HG22	1:A:126:ILE:CG1	2.30	0.61
1:B:26:TYR:CE2	1:B:112:LYS:HD3	2.35	0.61
1:B:122:THR:O	1:B:126:ILE:HG13	2.01	0.61
2:L:11:DC:H2''	2:L:12:DT:C5'	2.31	0.60
1:A:132:VAL:CG1	1:B:129:THR:HG21	2.24	0.59
1:A:8:ALA:O	1:A:12:VAL:HG23	2.03	0.59
1:A:84:SER:HB2	1:A:92:TYR:CE2	2.39	0.58
1:A:5:LEU:HD21	1:B:117:MET:CE	2.34	0.57
1:B:122:THR:HG23	1:B:125:GLU:OE2	2.05	0.56
2:L:7:DA:H2''	2:L:8:DG:H5'	1.86	0.56
1:A:121:ILE:CG2	1:A:126:ILE:CG1	2.84	0.55
3:M:21:DC:H2''	3:M:22:DT:H5''	1.88	0.55
2:L:16:DG:H2''	2:L:17:DA:H5'	1.89	0.55
1:B:140:MET:N	1:B:140:MET:SD	2.80	0.54
1:A:13:ASN:ND2	3:M:16:DG:H5'	2.23	0.54
1:A:62:SER:OG	2:L:17:DA:OP2	2.25	0.54
1:B:121:ILE:HG22	1:B:126:ILE:CG1	2.38	0.54
1:B:121:ILE:HG22	1:B:126:ILE:HG13	1.90	0.54
2:L:26:DA:H2''	2:L:27:DA:OP2	2.06	0.54
1:B:62:SER:OG	3:M:17:DA:OP2	2.26	0.53
1:A:74:GLU:OE1	1:A:81:ILE:HD12	2.08	0.52
1:B:24:LYS:HG3	1:B:29:GLU:HG3	1.91	0.52
1:B:26:TYR:HE2	1:B:112:LYS:HD3	1.74	0.52
1:A:53:LYS:HD2	1:A:53:LYS:O	2.10	0.52
1:A:129:THR:HG21	1:B:132:VAL:HG11	1.92	0.51
1:B:114:HIS:ND1	1:B:114:HIS:N	2.57	0.51
3:M:11:DA:H2''	3:M:12:DT:C5'	2.41	0.51
1:B:3:ASN:O	1:B:7:LYS:CG	2.54	0.50
1:B:23:ALA:HB1	1:B:28:VAL:HG23	1.92	0.50
1:A:121:ILE:CG2	1:A:126:ILE:HG13	2.41	0.50
1:B:13:ASN:ND2	2:L:16:DG:H5'	2.27	0.49
3:M:11:DA:H2''	3:M:12:DT:H5'	1.94	0.49
3:M:17:DA:H2''	3:M:18:DG:H5'	1.94	0.49
3:M:25:DT:H4'	3:M:25:DT:OP1	2.12	0.49
3:M:1:DT:H2''	3:M:2:DT:C5	2.48	0.48
1:B:134:ARG:O	1:B:137:ALA:HB3	2.14	0.48
1:B:130:LYS:HB3	1:B:134:ARG:NH1	2.28	0.48
1:A:123:LYS:O	1:A:127:ARG:N	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:LYS:HE3	1:A:112:LYS:HB2	1.66	0.47
1:A:37:HIS:CD2	1:B:6:GLN:HG3	2.49	0.47
3:M:16:DG:H2''	3:M:17:DA:H5'	1.97	0.47
1:A:5:LEU:HD21	1:B:117:MET:HE1	1.97	0.46
1:A:121:ILE:HG22	1:A:126:ILE:CD1	2.46	0.46
3:M:22:DT:H5''	3:M:22:DT:H6	1.81	0.45
2:L:11:DC:H2''	2:L:12:DT:O5'	2.17	0.45
1:A:127:ARG:O	1:A:130:LYS:HB2	2.17	0.45
1:B:136:LEU:O	1:B:139:ASN:HB2	2.17	0.45
1:A:121:ILE:HG22	1:A:126:ILE:HD11	1.99	0.44
1:A:117:MET:O	1:B:139:ASN:ND2	2.49	0.44
1:A:26:TYR:OH	1:A:112:LYS:HD2	2.17	0.44
1:A:113:LEU:CD2	1:B:5:LEU:HD11	2.48	0.43
2:L:18:DT:H2''	2:L:19:DA:C8	2.53	0.43
1:A:121:ILE:CG2	1:A:126:ILE:HD11	2.48	0.43
1:B:37:HIS:ND1	1:B:59:LEU:HD23	2.33	0.43
3:M:11:DA:H2''	3:M:12:DT:O5'	2.19	0.43
1:A:113:LEU:HD21	1:B:5:LEU:HD21	2.01	0.43
3:M:22:DT:H5''	3:M:22:DT:C6	2.54	0.43
1:A:140:MET:CG	1:B:18:TYR:CD2	3.02	0.42
2:L:18:DT:H1'	2:L:19:DA:H5'	2.01	0.42
1:A:31:LEU:O	1:A:36:GLY:HA3	2.20	0.42
2:L:15:DC:C2	2:L:16:DG:C8	3.08	0.42
1:A:6:GLN:CD	1:B:37:HIS:HB2	2.39	0.42
1:A:63:LYS:HB2	1:A:63:LYS:HE2	1.66	0.41
1:B:6:GLN:HA	1:B:6:GLN:OE1	2.20	0.41
2:L:18:DT:H2''	2:L:19:DA:H8	1.84	0.41
1:A:37:HIS:HB2	1:B:6:GLN:CD	2.41	0.41
2:L:26:DA:H1'	2:L:27:DA:H5'	2.00	0.41
1:A:4:PRO:HD2	1:B:114:HIS:HE2	1.85	0.41
2:L:23:DA:H2''	2:L:24:DT:O4'	2.21	0.41
1:B:95:LEU:HD23	1:B:95:LEU:HA	1.73	0.41
1:A:140:MET:HG3	1:B:18:TYR:HD2	1.83	0.41
1:B:3:ASN:HA	1:B:4:PRO:HD3	1.99	0.41
1:A:9:ARG:HH12	1:B:34:PRO:HD3	1.86	0.40
1:B:31:LEU:O	1:B:36:GLY:HA3	2.21	0.40
1:B:38:LEU:HD12	1:B:38:LEU:HA	1.85	0.40
3:M:7:DA:C2'	3:M:8:DG:H5'	2.50	0.40
1:B:113:LEU:HA	1:B:113:LEU:HD12	1.80	0.40
1:A:89:ARG:HD3	2:L:24:DT:O2	2.21	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	137/153 (90%)	132 (96%)	5 (4%)	0	100	100
1	B	138/153 (90%)	131 (95%)	7 (5%)	0	100	100
All	All	275/306 (90%)	263 (96%)	12 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	124/138 (90%)	110 (89%)	14 (11%)	6	11
1	B	125/138 (91%)	116 (93%)	9 (7%)	14	28
All	All	249/276 (90%)	226 (91%)	23 (9%)	9	18

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	6	GLN
1	A	49	ASP
1	A	53	LYS
1	A	62	SER
1	A	63	LYS
1	A	75	LYS
1	A	113	LEU

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Mol	Chain	Res	Type
1	A	115	SER
1	A	117	MET
1	A	122	THR
1	A	132	VAL
1	A	135	THR
1	A	140	MET
1	B	3	ASN
1	B	15	LEU
1	B	22	TYR
1	B	53	LYS
1	B	62	SER
1	B	75	LYS
1	B	84	SER
1	B	132	VAL
1	B	140	MET

Sometimes 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](#)

There are no ligands in this entry.

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, 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	139/153 (90%)	0.54	12 (8%) 10 10	39, 82, 150, 166	0
1	B	140/153 (91%)	0.54	15 (10%) 6 5	40, 86, 154, 194	0
2	L	28/28 (100%)	-0.59	0 100 100	42, 73, 114, 133	0
3	M	28/28 (100%)	-0.66	0 100 100	43, 74, 108, 122	0
All	All	335/362 (92%)	0.34	27 (8%) 12 12	39, 81, 150, 194	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	LEU	4.9
1	A	117	MET	4.8
1	B	11	LEU	4.0
1	B	5	LEU	3.9
1	A	18	TYR	3.8
1	B	15	LEU	3.6
1	B	117	MET	3.5
1	B	6	GLN	3.4
1	A	15	LEU	3.1
1	B	22	TYR	3.1
1	A	127	ARG	3.1
1	A	26	TYR	3.1
1	B	131	LYS	2.8
1	B	123	LYS	2.8
1	B	9	ARG	2.8
1	A	48	LYS	2.8
1	B	126	ILE	2.6
1	B	109	PHE	2.5
1	A	133	ILE	2.5
1	B	132	VAL	2.4
1	A	134	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	21	ARG	2.4
1	A	137	ALA	2.3
1	B	8	ALA	2.2
1	B	133	ILE	2.1
1	A	22	TYR	2.0
1	A	113	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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.