



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:46 PM UTC

PDB ID : 3LRR / pdb_00003lrr
Title : Crystal structure of human RIG-I CTD bound to a 12 bp AU rich 5' ppp dsRNA
Authors : Li, P.
Deposited on : 2010-02-11
Resolution : 2.15 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

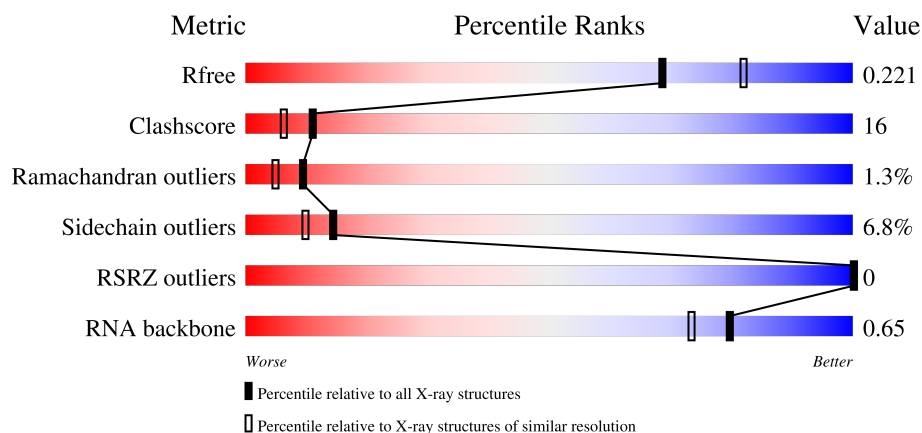
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)
RNA backbone	3983	1112 (2.50-1.82)

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	121	67% 25% 8%
1	B	121	52% 40% 6%
2	C	12	50% 33% 17%
2	D	12	50% 33% 17%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2522 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 Probable ATP-dependent RNA helicase DDX58.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	121	Total	C	N	O	S	0	0	0
			999	648	169	175	7			
1	B	121	Total	C	N	O	S	0	0	0
			999	648	169	175	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	829	SER	CYS	engineered mutation	UNP O95786
B	829	SER	CYS	engineered mutation	UNP O95786

- Molecule 2 is a RNA chain called RNA (5'-R(*(ATP)P*UP*AP*UP*AP*UP*AP*UP*AP*UP*AP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	P	0	0	0
			261	114	42	91	14			
2	D	12	Total	C	N	O	P	0	0	0
			261	114	42	91	14			

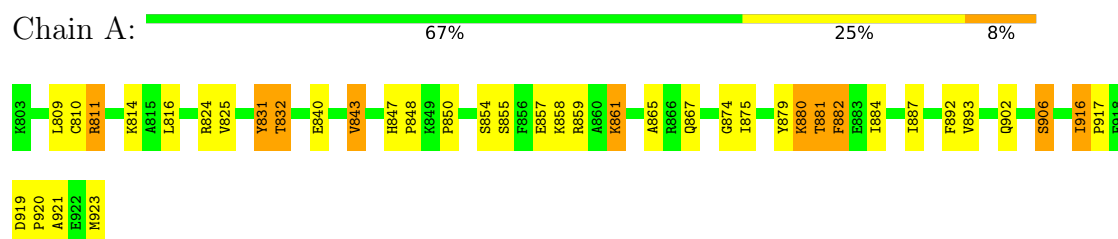
- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		

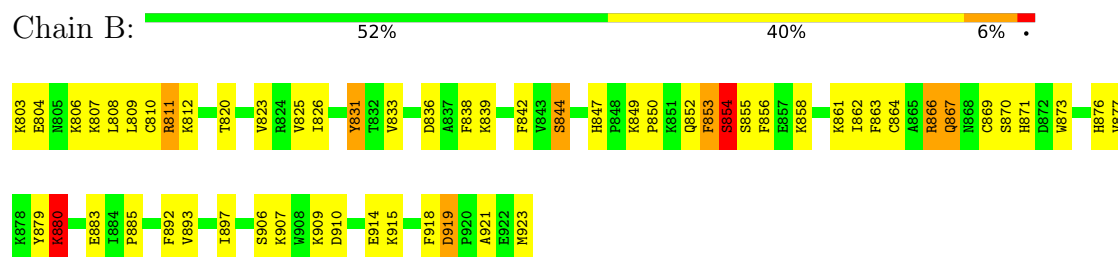
3 Residue-property plots

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: Probable ATP-dependent RNA helicase DDX58



- Molecule 1: Probable ATP-dependent RNA helicase DDX58



- Molecule 2: RNA (5'-R*(ATP)P*UP*AP*UP*AP*UP*AP*UP*AP*UP*AP*U)-3')



- Molecule 2: RNA (5'-R*(ATP)P*UP*AP*UP*AP*UP*AP*UP*AP*UP*AP*U)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	83.01Å 83.01Å 111.22Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.19 – 2.15 30.19 – 2.15	Depositor EDS
% Data completeness (in resolution range)	91.8 (30.19-2.15) 91.8 (30.19-2.15)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.24 (at 2.16Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.187 , 0.228 0.190 , 0.221	Depositor DCC
R_{free} test set	1452 reflections (6.59%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 23.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.480 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2522	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.42% 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: ATP, ZN

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	0.59	0/1027	1.10	7/1380 (0.5%)
1	B	0.59	0/1027	1.16	14/1380 (1.0%)
2	C	0.35	0/256	0.67	0/395
2	D	0.34	0/256	0.64	0/395
All	All	0.55	0/2566	1.04	21/3550 (0.6%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	853	PHE	CA-C-N	-7.26	110.74	120.54
1	B	853	PHE	C-N-CA	-7.26	110.74	120.54
1	A	882	PHE	N-CA-C	6.99	121.24	110.20
1	A	906	SER	N-CA-C	-6.35	106.07	113.88
1	A	884	ILE	N-CA-C	6.19	115.88	108.45
1	B	831	TYR	N-CA-C	6.18	118.98	108.90
1	B	880	LYS	CB-CA-C	-6.05	109.00	117.23
1	B	919	ASP	CA-C-N	5.90	125.77	119.28
1	B	919	ASP	C-N-CA	5.90	125.77	119.28
1	A	887	ILE	N-CA-C	5.75	117.89	108.97
1	A	916	ILE	N-CA-C	5.62	114.14	107.73
1	B	897	ILE	N-CA-C	5.58	116.31	110.62
1	A	831	TYR	N-CA-C	5.57	117.98	108.90
1	B	892	PHE	N-CA-C	5.53	118.58	109.85
1	B	854	SER	N-CA-CB	5.37	117.94	109.94
1	B	866	ARG	N-CA-C	5.36	117.95	109.96
1	A	881	THR	N-CA-C	-5.35	106.14	113.30
1	B	854	SER	N-CA-C	5.22	117.38	111.11
1	B	854	SER	CA-C-N	-5.03	114.14	121.99
1	B	854	SER	C-N-CA	-5.03	114.14	121.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	906	SER	N-CA-C	-5.03	107.69	113.88

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	999	0	993	27	0
1	B	999	0	993	44	0
2	C	261	0	127	8	0
2	D	261	0	127	6	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
All	All	2522	0	2240	78	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 16.

All (78) 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:B:812:LYS:HD3	1:B:869:CYS:SG	2.20	0.81
1:A:847:HIS:NE2	2:C:1:ATP:O2B	2.18	0.74
1:B:854:SER:HB2	2:C:12:U:O2'	1.87	0.73
1:B:847:HIS:CD2	1:B:861:LYS:HG2	2.27	0.69
1:B:803:LYS:N	1:B:914:GLU:HB2	2.07	0.68
1:B:864:CYS:O	1:B:870:SER:HA	1.96	0.66
1:A:847:HIS:NE2	1:A:858:LYS:HE3	2.11	0.66
1:B:811:ARG:HA	1:B:893:VAL:HG23	1.79	0.65
1:B:907:LYS:HB3	1:B:909:LYS:HG2	1.78	0.64
1:B:825:VAL:HG21	1:B:915:LYS:HB3	1.79	0.64
1:A:861:LYS:HE2	1:A:874:GLY:O	1.99	0.62
1:B:807:LYS:HE2	1:B:809:LEU:HD21	1.80	0.62
2:D:9:A:H2'	2:D:10:U:C6	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:9:A:H2'	2:D:10:U:H6	1.66	0.60
1:A:847:HIS:HE2	1:A:858:LYS:HE3	1.65	0.60
1:B:804:GLU:HB2	1:B:806:LYS:NZ	2.16	0.60
1:A:843:VAL:HG22	1:A:865:ALA:HA	1.82	0.60
1:B:836:ASP:HA	1:B:839:LYS:HD2	1.85	0.59
1:A:843:VAL:CG2	1:A:865:ALA:HA	2.33	0.58
1:B:847:HIS:O	1:B:850:PRO:HD3	2.04	0.57
1:B:909:LYS:NZ	1:B:910:ASP:OD1	2.35	0.56
1:B:842:PHE:HB2	1:B:863:PHE:O	2.06	0.56
1:A:879:TYR:O	1:A:880:LYS:C	2.48	0.55
1:B:858:LYS:HA	1:B:877:VAL:HG12	1.89	0.55
1:B:849:LYS:HE3	2:D:1:ATP:O1G	2.07	0.55
1:B:853:PHE:O	1:B:854:SER:C	2.49	0.54
1:B:867:GLN:H	1:B:867:GLN:CD	2.16	0.53
1:B:918:PHE:HE1	1:B:923:MET:SD	2.30	0.53
1:A:811:ARG:HA	1:A:893:VAL:HG23	1.92	0.52
1:B:820:THR:HA	1:B:823:VAL:HG23	1.92	0.52
1:B:826:ILE:HG21	1:B:856:PHE:CZ	2.45	0.52
1:A:847:HIS:O	1:A:850:PRO:HD3	2.09	0.51
2:D:3:A:O2'	2:D:4:U:H5'	2.10	0.51
1:B:826:ILE:HG21	1:B:856:PHE:HZ	1.76	0.50
1:A:919:ASP:OD1	1:A:921:ALA:HB3	2.10	0.50
2:C:7:A:H2'	2:C:8:U:C6	2.45	0.50
1:A:879:TYR:O	1:A:882:PHE:N	2.45	0.50
1:B:919:ASP:OD1	1:B:921:ALA:HB3	2.13	0.48
1:B:855:SER:HB3	1:B:880:LYS:NZ	2.28	0.48
2:C:1:ATP:H2'	2:C:2:U:C6	2.49	0.48
1:A:858:LYS:HG3	1:A:875:ILE:HD12	1.96	0.48
1:A:825:VAL:HG22	1:A:831:TYR:CE2	2.49	0.47
1:A:919:ASP:OD1	1:A:921:ALA:N	2.47	0.47
1:B:810:CYS:HB2	1:B:873:TRP:HZ2	1.79	0.47
1:B:804:GLU:HB2	1:B:806:LYS:HZ3	1.78	0.47
1:B:918:PHE:CE1	1:B:923:MET:SD	3.08	0.47
1:B:854:SER:H	2:C:12:U:H2'	1.78	0.46
1:B:825:VAL:HG22	1:B:831:TYR:CE2	2.51	0.46
1:B:826:ILE:CG2	1:B:856:PHE:HZ	2.29	0.46
1:A:847:HIS:CD2	1:A:858:LYS:HG2	2.51	0.45
1:B:856:PHE:CD1	1:B:879:TYR:HB2	2.52	0.45
1:A:810:CYS:O	1:A:814:LYS:HA	2.16	0.45
1:B:864:CYS:HB3	1:B:871:HIS:H	1.82	0.44
1:A:824:ARG:HB2	1:A:832:THR:HG22	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:825:VAL:CG2	1:B:915:LYS:HB3	2.48	0.44
1:B:864:CYS:SG	1:B:866:ARG:HB2	2.57	0.44
1:B:907:LYS:HD3	1:B:909:LYS:HD3	2.00	0.44
1:B:844:SER:HB3	1:B:862:ILE:HG22	1.99	0.44
1:A:920:PRO:O	1:A:923:MET:N	2.52	0.43
1:A:861:LYS:NZ	2:C:1:ATP:O2A	2.44	0.43
1:B:833:VAL:HB	1:B:838:PHE:CD2	2.53	0.43
1:A:916:ILE:HA	1:A:917:PRO:HD3	1.96	0.42
1:A:857:GLU:O	1:A:859:ARG:HG3	2.19	0.42
1:A:809:LEU:O	1:A:892:PHE:HA	2.19	0.42
1:A:809:LEU:HD23	1:A:816:LEU:HA	2.01	0.42
2:C:6:U:O2'	2:C:7:A:H5'	2.20	0.42
1:A:919:ASP:HA	1:A:920:PRO:HD3	1.96	0.42
1:A:893:VAL:CG1	1:A:902:GLN:HB3	2.50	0.41
1:B:907:LYS:CD	1:B:909:LYS:HD3	2.50	0.41
1:B:847:HIS:NE2	2:D:1:ATP:O2B	2.51	0.41
1:B:880:LYS:HD2	1:B:880:LYS:HA	1.78	0.41
2:D:8:U:H2'	2:D:9:A:H8	1.85	0.41
1:A:847:HIS:HA	1:A:848:PRO:HD2	1.89	0.41
1:A:811:ARG:NH1	1:A:892:PHE:O	2.44	0.41
1:B:876:HIS:HE1	1:B:883:GLU:HB3	1.86	0.41
1:B:854:SER:N	2:C:12:U:H2'	2.36	0.40
1:B:876:HIS:HA	1:B:885:PRO:HA	2.04	0.40
1:B:826:ILE:HG12	1:B:918:PHE:CB	2.51	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	119/121 (98%)	110 (92%)	7 (6%)	2 (2%)	7 2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	119/121 (98%)	110 (92%)	8 (7%)	1 (1%)	16	10
All	All	238/242 (98%)	220 (92%)	15 (6%)	3 (1%)	9	5

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	867	GLN
1	A	880	LYS
1	B	867	GLN

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	110/110 (100%)	101 (92%)	9 (8%)	10	6
1	B	110/110 (100%)	104 (94%)	6 (6%)	19	15
All	All	220/220 (100%)	205 (93%)	15 (7%)	14	10

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	811	ARG
1	A	832	THR
1	A	840	GLU
1	A	843	VAL
1	A	854	SER
1	A	855	SER
1	A	861	LYS
1	A	881	THR
1	A	906	SER
1	B	808	LEU
1	B	811	ARG
1	B	844	SER
1	B	852	GLN

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Mol	Chain	Res	Type
1	B	854	SER
1	B	880	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	805	ASN
1	B	876	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	11/12 (91%)	1 (9%)	1 (9%)
2	D	11/12 (91%)	1 (9%)	1 (9%)
All	All	22/24 (91%)	2 (9%)	2 (9%)

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	8	U
2	D	8	U

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	C	1	ATP
2	D	1	ATP

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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

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 [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	121/121 (100%)	-1.36	0 100 100	20, 32, 60, 82	0
1	B	121/121 (100%)	-1.33	0 100 100	20, 34, 64, 87	0
2	C	11/12 (91%)	-1.39	0 100 100	31, 79, 101, 104	0
2	D	11/12 (91%)	-1.39	0 100 100	32, 76, 95, 96	0
All	All	264/266 (99%)	-1.35	0 100 100	20, 34, 80, 104	0

There are no RSRZ outliers to report.

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
3	ZN	A	1	1/1	1.00	0.01	33,33,33,33	0
3	ZN	B	2	1/1	1.00	0.01	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.