



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:26 PM UTC

PDB ID : 5EI9 / pdb_00005ei9
Title : Human PRDM9 allele-A ZnF Domain with Associated Recombination Hotspot DNA Sequence I
Authors : Patel, A.; Horton, J.R.; Wilson, G.G.; Zhang, X.; Cheng, X.
Deposited on : 2015-10-29
Resolution : 1.92 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
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

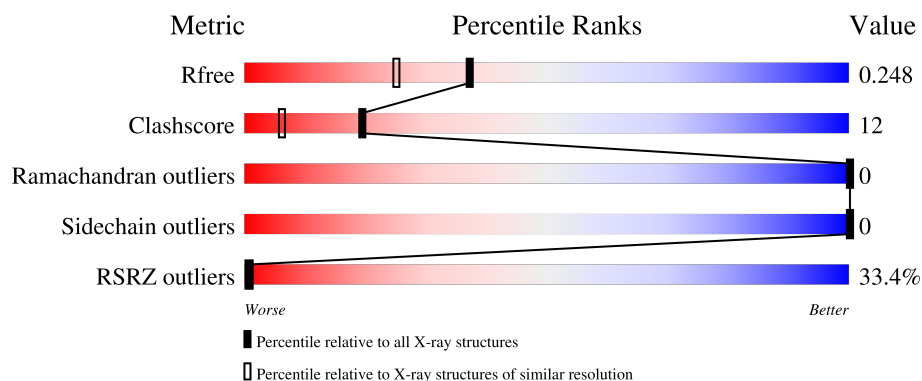
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	21	48% 67% 33%
1	C	21	38% 52% 48%
2	E	144	17% 65% 12% 23%
2	F	144	29% 60% 15% 24%
3	B	21	43% 76% 24%

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Mol	Chain	Length	Quality of chain
3	D	21	38% 71% 29%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3722 atoms, of which 16 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 DNA chain called DNA (5'-D(*AP*TP*CP*CP*AP*CP*GP*TP*GP*GP*CP*TP*AP*GP*GP*GP*AP*GP*GP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	21	Total	C	N	O	P	0	0	0
			431	204	84	123	20			
1	C	21	Total	C	N	O	P	0	0	0
			431	204	84	123	20			

- Molecule 2 is a protein called Histone-lysine N-methyltransferase PRDM9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	110	Total	C	H	N	O	S	0	0
			912	546	8	197	153	8		
2	E	111	Total	C	H	N	O	S	0	1
			925	554	8	201	154	8		0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	714	GLY	-	expression tag	UNP Q9NQV7
F	715	SER	-	expression tag	UNP Q9NQV7
E	714	GLY	-	expression tag	UNP Q9NQV7
E	715	SER	-	expression tag	UNP Q9NQV7

- Molecule 3 is a DNA chain called DNA (5'-D(*TP*GP*GP*CP*CP*TP*CP*CP*CP*TP*AP*GP*CP*CP*AP*CP*GP*TP*GP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	B	21	Total	C	N	O	P	0	0	0
			423	202	77	124	20			
3	D	21	Total	C	N	O	P	0	0	0
			424	202	77	125	20			

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	4	Total 4	Zn 4	0	0
4	E	4	Total 4	Zn 4	0	0

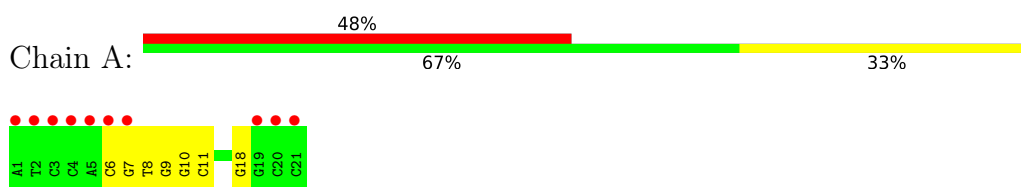
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	33	Total 33	O 33	0	0
5	F	31	Total 31	O 31	0	0
5	E	52	Total 52	O 52	0	0
5	C	18	Total 18	O 18	0	0
5	B	24	Total 24	O 24	0	0
5	D	10	Total 10	O 10	0	0

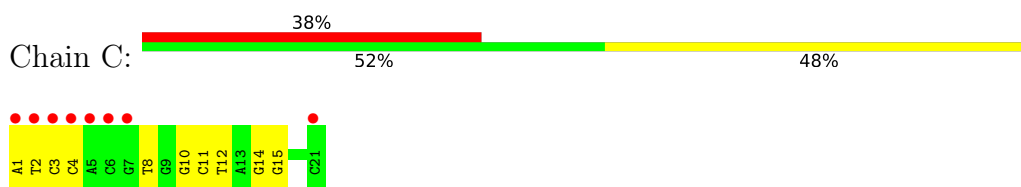
3 Residue-property plots [i](#)

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.

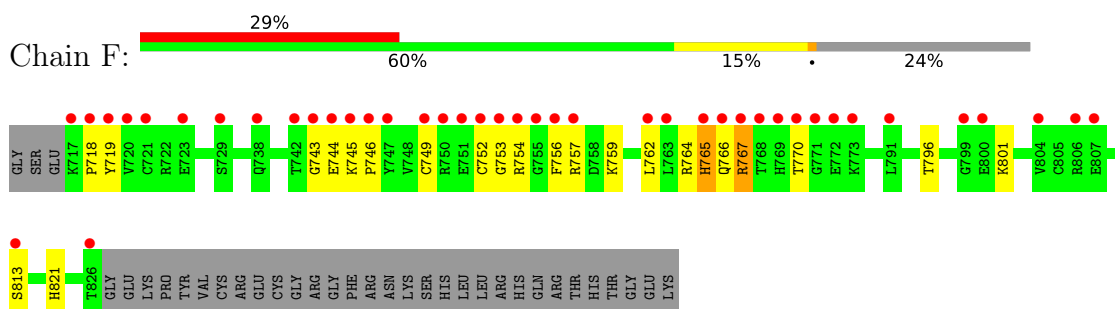
- Molecule 1: DNA (5'-D(*AP*TP*CP*CP*AP*CP*GP*TP*GP*GP*CP*TP*AP*GP*GP*GP*AP*GP*GP*CP*C)-3')



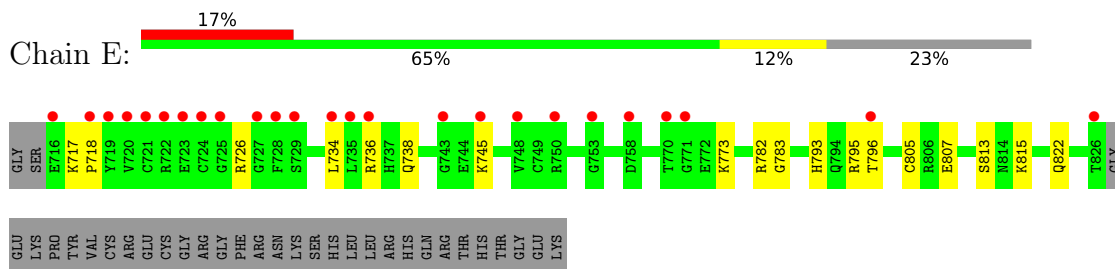
- Molecule 1: DNA (5'-D(*AP*TP*CP*CP*AP*CP*GP*TP*GP*GP*CP*TP*AP*GP*GP*GP*AP*GP*GP*CP*C)-3')



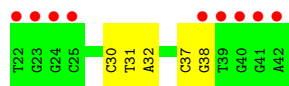
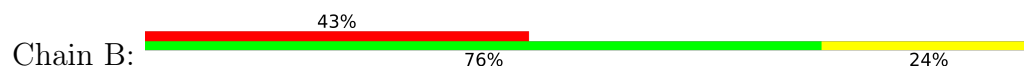
- Molecule 2: Histone-lysine N-methyltransferase PRDM9



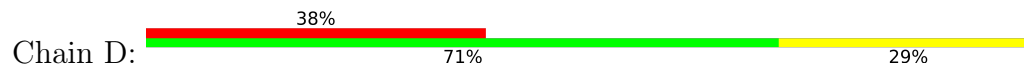
- Molecule 2: Histone-lysine N-methyltransferase PRDM9



- Molecule 3: DNA (5'-D(*TP*GP*GP*CP*CP*TP*CP*CP*CP*TP*AP*GP*CP*CP*AP*CP*GP*TP*GP*GP*A)-3')



- Molecule 3: DNA (5'-D(*TP*GP*GP*CP*CP*TP*CP*CP*CP*TP*AP*GP*CP*CP*AP*CP*GP*TP*GP*GP*A)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	44.33Å 55.74Å 76.04Å 89.85° 75.50° 86.58°	Depositor
Resolution (Å)	32.98 – 1.92 32.98 – 1.92	Depositor EDS
% Data completeness (in resolution range)	97.5 (32.98-1.92) 88.9 (32.98-1.92)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.32 (at 1.92Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.214 , 0.237 0.227 , 0.248	Depositor DCC
R_{free} test set	1995 reflections (3.79%)	wwPDB-VP
Wilson B-factor (Å ²)	36.3	Xtriage
Anisotropy	0.620	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,-h+l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3722	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.54 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.5031e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: 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	0.47	0/484	0.71	0/746
1	C	0.41	0/484	0.56	0/746
2	E	0.62	0/942	0.89	2/1256 (0.2%)
2	F	0.54	0/926	0.93	5/1235 (0.4%)
3	B	0.43	0/473	0.55	0/728
3	D	0.33	0/474	0.49	0/729
All	All	0.50	0/3783	0.75	7/5440 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	745	LYS	CA-C-N	6.86	127.00	119.87
2	F	745	LYS	C-N-CA	6.86	127.00	119.87
2	F	753	GLY	N-CA-C	-6.58	105.43	114.64
2	E	745	LYS	CA-C-N	6.47	126.17	119.64
2	E	745	LYS	C-N-CA	6.47	126.17	119.64
2	F	767	ARG	N-CA-C	-6.03	105.00	112.90
2	F	765	HIS	N-CA-C	-5.41	103.35	110.43

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	431	0	236	15	0
1	C	431	0	236	17	0
2	E	917	8	880	18	0
2	F	904	8	865	21	0
3	B	423	0	234	6	0
3	D	424	0	237	4	0
4	E	4	0	0	0	0
4	F	4	0	0	0	0
5	A	33	0	0	1	0
5	B	24	0	0	0	0
5	C	18	0	0	0	0
5	D	10	0	0	0	0
5	E	52	0	0	2	0
5	F	31	0	0	0	0
All	All	3706	16	2688	70	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 12.

All (70) 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:C:14:DG:H2''	1:C:15:DG:H5'	1.43	1.01
1:C:14:DG:H5''	1:C:14:DG:H8	1.33	0.92
1:C:14:DG:H5''	1:C:14:DG:C8	2.11	0.86
2:E:783:GLY:O	5:E:1101:HOH:O	2.01	0.79
1:A:11:DC:OP1	2:E:782:ARG:NH2	2.17	0.78
1:A:6:DC:H2'	1:A:7:DG:C8	2.21	0.74
1:C:14:DG:H2''	1:C:15:DG:C5'	2.21	0.69
2:E:734:LEU:O	2:E:738:GLN:HG2	1.94	0.67
1:A:18:DG:N7	2:E:736:ARG:NH1	2.41	0.67
2:F:749:CYS:HB3	2:F:752:CYS:HB2	1.78	0.66
1:A:18:DG:O6	2:E:736:ARG:NH2	2.21	0.66
2:E:717:LYS:CB	2:E:718:PRO:HD2	2.28	0.63
2:F:765:HIS:O	2:F:766:GLN:HB2	1.99	0.61
2:E:815:LYS:HE2	3:B:31:DT:OP2	2.02	0.59
2:F:821:HIS:ND1	1:C:8:DT:OP1	2.35	0.59
1:A:10:DG:H3'	2:E:796:THR:HG21	1.86	0.58
2:F:756:PHE:HE2	1:C:14:DG:H5'	1.68	0.58
3:D:41:DG:H2'	3:D:42:DA:O4'	2.04	0.57
1:A:9:DG:H4'	1:A:10:DG:OP1	2.05	0.57
1:A:6:DC:H2''	1:A:7:DG:O5'	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:822:GLN:NE2	5:E:1103:HOH:O	2.38	0.55
2:F:752:CYS:SG	2:F:765:HIS:NE2	2.80	0.55
1:C:3:DC:H2''	1:C:4:DC:H5'	1.89	0.54
3:D:30:DC:H2''	3:D:31:DT:H5'	1.89	0.54
1:A:6:DC:H2'	1:A:7:DG:H8	1.73	0.53
3:D:39:DT:H2''	3:D:40:DG:C8	2.43	0.53
5:A:107:HOH:O	2:E:773:LYS:HE2	2.08	0.53
1:A:11:DC:OP2	2:E:796:THR:HG21	2.10	0.52
3:B:30:DC:H2''	3:B:31:DT:H5'	1.90	0.52
1:C:14:DG:H2'	1:C:15:DG:C8	2.45	0.52
1:C:14:DG:C2'	1:C:15:DG:H5'	2.30	0.51
2:F:752:CYS:CB	2:F:754:ARG:H	2.24	0.50
2:E:717:LYS:CB	2:E:718:PRO:CD	2.90	0.50
1:A:6:DC:H2''	1:A:7:DG:C5'	2.42	0.49
2:F:744:GLU:OE2	2:F:746:PRO:HG3	2.12	0.49
1:A:6:DC:H4'	1:A:7:DG:OP1	2.13	0.49
2:F:765:HIS:CD2	2:F:766:GLN:N	2.81	0.48
2:F:765:HIS:CD2	2:F:766:GLN:H	2.31	0.48
2:F:752:CYS:HB3	2:F:754:ARG:HG3	1.95	0.48
2:F:743:GLY:HA2	2:F:757:ARG:HH21	1.78	0.48
1:C:14:DG:H2'	1:C:15:DG:H8	1.77	0.48
1:A:9:DG:H2''	1:A:10:DG:O5'	2.14	0.47
3:B:30:DC:C2'	3:B:31:DT:H5'	2.45	0.47
2:F:765:HIS:ND1	1:C:14:DG:OP1	2.48	0.47
2:F:767:ARG:HA	2:F:770:THR:OG1	2.15	0.47
3:B:37:DC:H2''	3:B:38:DG:C8	2.50	0.46
1:C:1:DA:H4'	1:C:2:DT:O5'	2.16	0.46
1:C:11:DC:H2''	1:C:12:DT:C5'	2.46	0.45
2:F:801:LYS:HA	2:F:813:SER:HA	1.97	0.45
2:E:726[B]:ARG:HG3	2:E:726[B]:ARG:NH1	2.32	0.45
1:A:7:DG:C2'	1:A:8:DT:H71	2.46	0.45
2:F:752:CYS:HB3	2:F:754:ARG:H	1.82	0.45
3:D:40:DG:H2'	3:D:41:DG:C8	2.52	0.44
2:F:796:THR:HG21	1:C:10:DG:H3'	1.99	0.44
2:E:795:ARG:HD2	2:E:813:SER:O	2.17	0.43
3:B:31:DT:H2''	3:B:32:DA:C8	2.53	0.43
2:E:726[B]:ARG:HG3	2:E:726[B]:ARG:HH11	1.83	0.43
1:A:7:DG:H2'	1:A:8:DT:H71	2.00	0.43
1:C:14:DG:C2'	1:C:15:DG:C5'	2.93	0.42
2:F:759:LYS:O	2:F:762:LEU:HB3	2.19	0.42
2:F:765:HIS:O	2:F:766:GLN:CB	2.65	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:805:CYS:SG	2:E:807:GLU:HB3	2.58	0.42
2:F:718:PRO:HG2	2:F:719:TYR:CD2	2.54	0.42
2:F:764:ARG:O	2:F:767:ARG:HB2	2.21	0.41
1:A:10:DG:H3'	2:E:796:THR:CG2	2.50	0.41
2:E:793:HIS:O	2:E:796:THR:HG23	2.21	0.41
2:F:765:HIS:C	2:F:767:ARG:H	2.28	0.40
3:B:37:DC:H2''	3:B:38:DG:H8	1.85	0.40
1:C:2:DT:H2''	1:C:3:DC:O5'	2.20	0.40
1:C:11:DC:H2''	1:C:12:DT:H5'	2.03	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	
2	E	110/144 (76%)	108 (98%)	2 (2%)	0	100	100
2	F	108/144 (75%)	105 (97%)	3 (3%)	0	100	100
All	All	218/288 (76%)	213 (98%)	5 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	
2	E	98/128 (77%)	98 (100%)	0	100	100
2	F	97/128 (76%)	97 (100%)	0	100	100
All	All	195/256 (76%)	195 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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 8 ligands modelled in this entry, 8 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

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	21/21 (100%)	2.08	10 (47%) 0 0	34, 51, 83, 87	0
1	C	21/21 (100%)	1.95	8 (38%) 1 0	42, 54, 99, 100	0
2	E	111/144 (77%)	1.51	25 (22%) 2 2	30, 43, 71, 78	1 (0%)
2	F	110/144 (76%)	1.91	42 (38%) 1 0	35, 50, 75, 82	0
3	B	21/21 (100%)	2.12	9 (42%) 0 0	38, 55, 85, 88	0
3	D	21/21 (100%)	2.29	8 (38%) 1 0	46, 65, 97, 101	0
All	All	305/372 (81%)	1.82	102 (33%) 1 1	30, 49, 83, 101	1 (0%)

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	DA	6.3
3	B	22	DT	5.8
3	D	22	DT	5.0
2	F	752	CYS	4.9
1	C	1	DA	4.6
2	E	826	THR	4.6
3	B	23	DG	4.5
3	D	40	DG	4.4
2	F	770	THR	4.4
2	F	768	THR	4.4
2	E	723	GLU	4.3
2	E	721	CYS	4.2
3	B	42	DA	4.2
2	F	771	GLY	4.1
1	A	21	DC	4.0
3	D	39	DT	4.0
2	E	720	VAL	4.0
3	B	40	DG	4.0
2	F	717	LYS	3.9

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Mol	Chain	Res	Type	RSRZ
2	F	753	GLY	3.8
1	A	2	DT	3.7
3	D	38	DG	3.7
3	D	41	DG	3.7
1	C	2	DT	3.7
1	C	3	DC	3.7
3	B	41	DG	3.7
3	D	42	DA	3.6
2	F	729	SER	3.5
2	F	750	ARG	3.5
2	F	765	HIS	3.5
2	E	734	LEU	3.4
1	A	6	DC	3.4
2	E	716	GLU	3.3
2	F	813	SER	3.3
2	F	766	GLN	3.2
1	A	4	DC	3.2
2	F	723	GLU	3.2
2	F	749	CYS	3.2
1	C	5	DA	3.2
2	F	747	TYR	3.1
2	E	728	PHE	3.1
2	F	769	HIS	3.1
3	B	24	DG	3.1
2	F	756	PHE	3.1
2	F	751	GLU	3.1
1	C	7	DG	3.1
1	A	5	DA	3.0
1	C	6	DC	3.0
2	F	791	LEU	3.0
2	F	800	GLU	3.0
2	F	804	VAL	3.0
2	F	746	PRO	3.0
2	F	754	ARG	3.0
2	F	762	LEU	2.9
3	D	23	DG	2.9
2	F	826	THR	2.9
1	A	7	DG	2.9
2	E	718	PRO	2.8
2	E	735	LEU	2.8
2	E	727	GLY	2.8
1	A	3	DC	2.8

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Mol	Chain	Res	Type	RSRZ
2	E	729	SER	2.8
2	F	757	ARG	2.7
3	B	38	DG	2.7
2	F	720	VAL	2.6
2	E	758	ASP	2.6
1	C	21	DC	2.6
2	F	719	TYR	2.5
2	E	724	CYS	2.5
2	F	738	GLN	2.5
2	E	753	GLY	2.5
2	E	750	ARG	2.5
3	D	25	DC	2.5
3	B	39	DT	2.5
2	E	725	GLY	2.5
1	A	19	DG	2.5
3	B	25	DC	2.4
2	F	743	GLY	2.4
1	C	4	DC	2.4
2	F	767	ARG	2.4
2	F	745	LYS	2.4
2	F	763	LEU	2.4
2	E	722	ARG	2.4
2	E	770	THR	2.4
2	F	742	THR	2.3
1	A	20	DC	2.3
2	F	772	GLU	2.3
2	E	745	LYS	2.3
2	F	755	GLY	2.2
2	F	773	LYS	2.2
2	F	806	ARG	2.2
2	F	718	PRO	2.2
2	F	807	GLU	2.1
2	E	719	TYR	2.1
2	F	744	GLU	2.1
2	E	771	GLY	2.1
2	F	721	CYS	2.1
2	E	796	THR	2.0
2	E	743	GLY	2.0
2	E	748	VAL	2.0
2	E	736	ARG	2.0
2	F	799	GLY	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
4	ZN	F	1002	1/1	0.93	0.09	76,76,76,76	0
4	ZN	F	1004	1/1	0.96	0.06	54,54,54,54	0
4	ZN	E	1003	1/1	0.96	0.05	37,37,37,37	0
4	ZN	E	1004	1/1	0.96	0.06	42,42,42,42	0
4	ZN	F	1003	1/1	0.97	0.05	41,41,41,41	0
4	ZN	E	1001	1/1	0.97	0.04	49,49,49,49	0
4	ZN	F	1001	1/1	0.98	0.04	38,38,38,38	0
4	ZN	E	1002	1/1	0.98	0.04	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.