



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:59 PM UTC

PDB ID : 5K1S / pdb_00005k1s
Title : crystal structure of AibC
Authors : Bock, T.; Mueller, R.; Blankenfeldt, W.
Deposited on : 2016-05-18
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
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

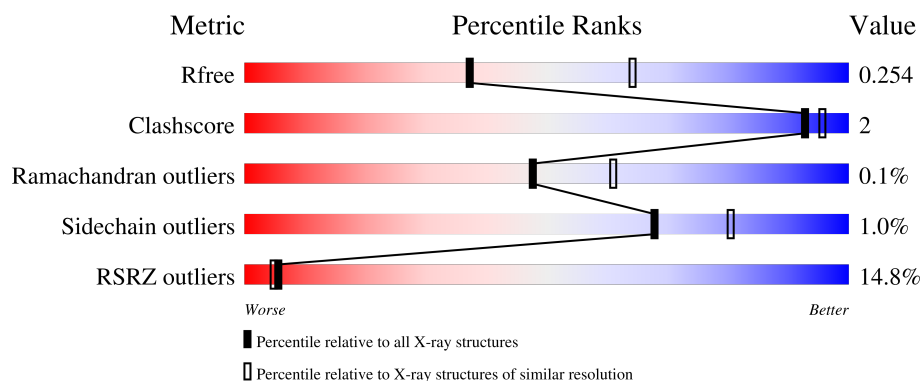
1 Overall quality at a glance

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	362	4% 93% 6%
1	B	362	24% 87% 5% 8%
1	C	362	11% 94% 5%
1	D	362	17% 90% 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ZN	C	402	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 19587 atoms, of which 9631 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 Oxidoreductase, zinc-binding dehydrogenase family.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	358	Total	C	H	N	O	S	0	1	0
			5116	1607	2529	476	488	16			
1	C	358	Total	C	H	N	O	S	0	0	0
			5033	1588	2480	469	480	16			
1	B	333	Total	C	H	N	O	S	0	0	0
			4641	1453	2298	426	449	15			
1	D	333	Total	C	H	N	O	S	0	0	0
			4682	1467	2324	423	452	16			

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	MET	-	initiating methionine	UNP Q1D4I2
A	-15	GLY	-	expression tag	UNP Q1D4I2
A	-14	HIS	-	expression tag	UNP Q1D4I2
A	-13	HIS	-	expression tag	UNP Q1D4I2
A	-12	HIS	-	expression tag	UNP Q1D4I2
A	-11	HIS	-	expression tag	UNP Q1D4I2
A	-10	HIS	-	expression tag	UNP Q1D4I2
A	-9	HIS	-	expression tag	UNP Q1D4I2
A	-8	ALA	-	expression tag	UNP Q1D4I2
A	-7	GLU	-	expression tag	UNP Q1D4I2
A	-6	ASN	-	expression tag	UNP Q1D4I2
A	-5	LEU	-	expression tag	UNP Q1D4I2
A	-4	TYR	-	expression tag	UNP Q1D4I2
A	-3	PHE	-	expression tag	UNP Q1D4I2
A	-2	GLN	-	expression tag	UNP Q1D4I2
A	-1	GLY	-	expression tag	UNP Q1D4I2
A	0	HIS	-	expression tag	UNP Q1D4I2
C	-16	MET	-	initiating methionine	UNP Q1D4I2
C	-15	GLY	-	expression tag	UNP Q1D4I2
C	-14	HIS	-	expression tag	UNP Q1D4I2
C	-13	HIS	-	expression tag	UNP Q1D4I2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-12	HIS	-	expression tag	UNP Q1D4I2
C	-11	HIS	-	expression tag	UNP Q1D4I2
C	-10	HIS	-	expression tag	UNP Q1D4I2
C	-9	HIS	-	expression tag	UNP Q1D4I2
C	-8	ALA	-	expression tag	UNP Q1D4I2
C	-7	GLU	-	expression tag	UNP Q1D4I2
C	-6	ASN	-	expression tag	UNP Q1D4I2
C	-5	LEU	-	expression tag	UNP Q1D4I2
C	-4	TYR	-	expression tag	UNP Q1D4I2
C	-3	PHE	-	expression tag	UNP Q1D4I2
C	-2	GLN	-	expression tag	UNP Q1D4I2
C	-1	GLY	-	expression tag	UNP Q1D4I2
C	0	HIS	-	expression tag	UNP Q1D4I2
B	-16	MET	-	initiating methionine	UNP Q1D4I2
B	-15	GLY	-	expression tag	UNP Q1D4I2
B	-14	HIS	-	expression tag	UNP Q1D4I2
B	-13	HIS	-	expression tag	UNP Q1D4I2
B	-12	HIS	-	expression tag	UNP Q1D4I2
B	-11	HIS	-	expression tag	UNP Q1D4I2
B	-10	HIS	-	expression tag	UNP Q1D4I2
B	-9	HIS	-	expression tag	UNP Q1D4I2
B	-8	ALA	-	expression tag	UNP Q1D4I2
B	-7	GLU	-	expression tag	UNP Q1D4I2
B	-6	ASN	-	expression tag	UNP Q1D4I2
B	-5	LEU	-	expression tag	UNP Q1D4I2
B	-4	TYR	-	expression tag	UNP Q1D4I2
B	-3	PHE	-	expression tag	UNP Q1D4I2
B	-2	GLN	-	expression tag	UNP Q1D4I2
B	-1	GLY	-	expression tag	UNP Q1D4I2
B	0	HIS	-	expression tag	UNP Q1D4I2
D	-16	MET	-	initiating methionine	UNP Q1D4I2
D	-15	GLY	-	expression tag	UNP Q1D4I2
D	-14	HIS	-	expression tag	UNP Q1D4I2
D	-13	HIS	-	expression tag	UNP Q1D4I2
D	-12	HIS	-	expression tag	UNP Q1D4I2
D	-11	HIS	-	expression tag	UNP Q1D4I2
D	-10	HIS	-	expression tag	UNP Q1D4I2
D	-9	HIS	-	expression tag	UNP Q1D4I2
D	-8	ALA	-	expression tag	UNP Q1D4I2
D	-7	GLU	-	expression tag	UNP Q1D4I2
D	-6	ASN	-	expression tag	UNP Q1D4I2
D	-5	LEU	-	expression tag	UNP Q1D4I2

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-4	TYR	-	expression tag	UNP Q1D4I2
D	-3	PHE	-	expression tag	UNP Q1D4I2
D	-2	GLN	-	expression tag	UNP Q1D4I2
D	-1	GLY	-	expression tag	UNP Q1D4I2
D	0	HIS	-	expression tag	UNP Q1D4I2

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		
2	C	2	Total	Zn	0	0
			2	2		
2	B	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		

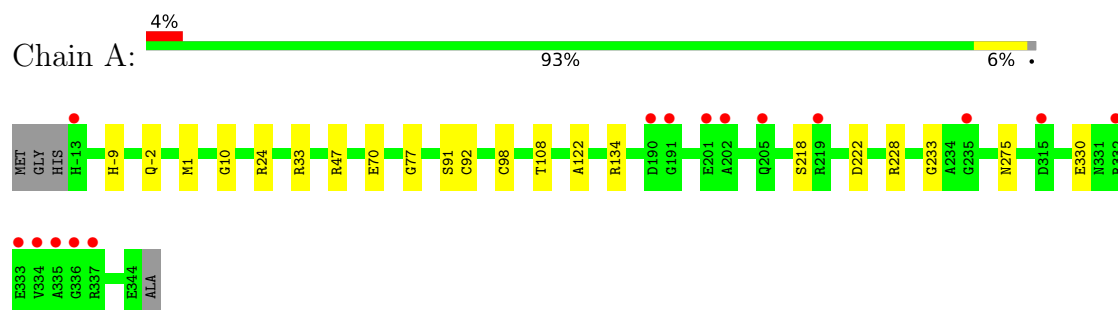
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	42	Total	O	0	0
			42	42		
3	C	35	Total	O	0	0
			35	35		
3	B	18	Total	O	0	0
			18	18		
3	D	12	Total	O	0	0
			12	12		

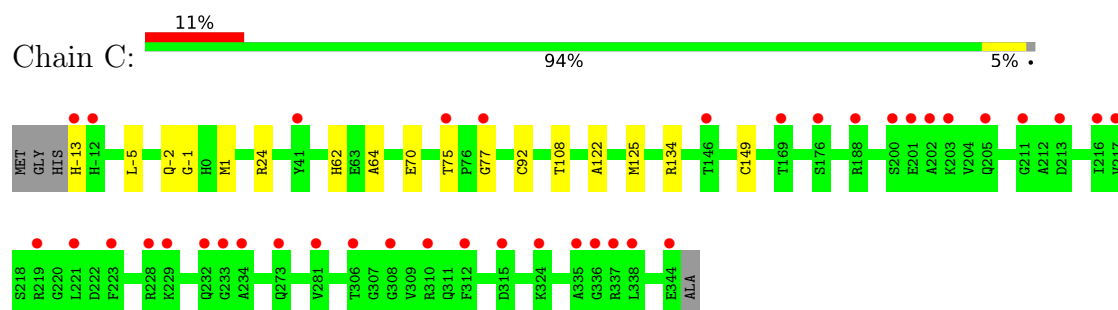
3 Residue-property plots [i](#)

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.

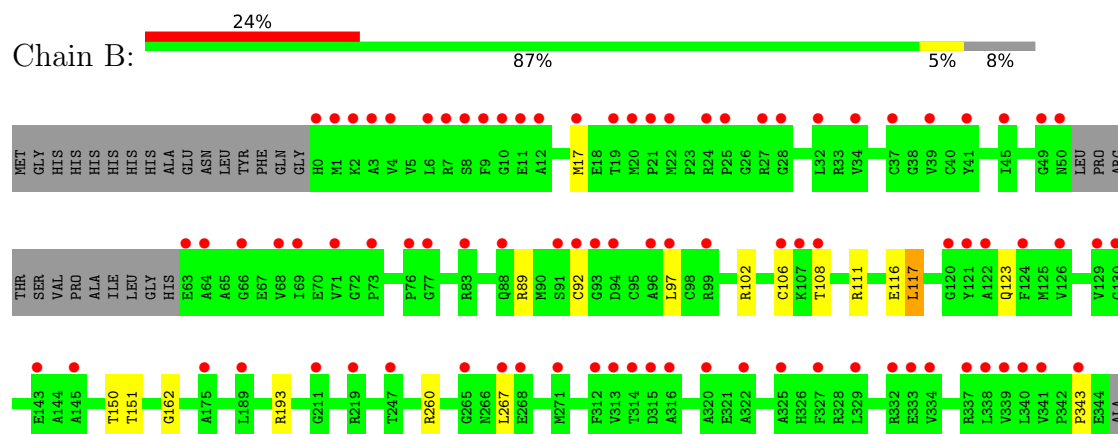
- Molecule 1: Oxidoreductase, zinc-binding dehydrogenase family



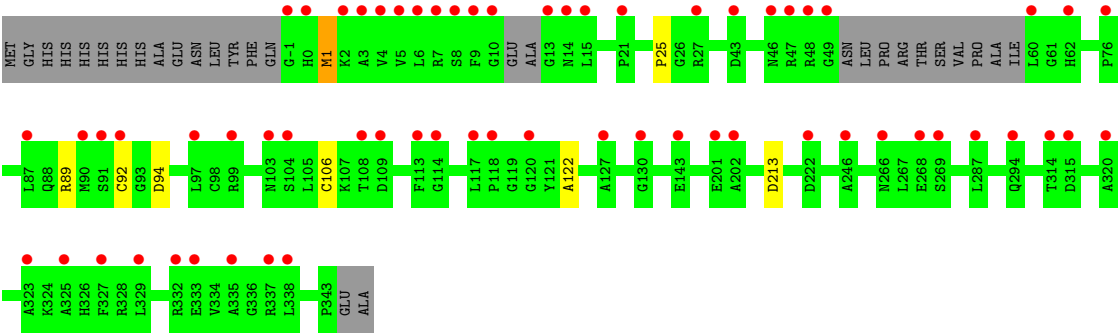
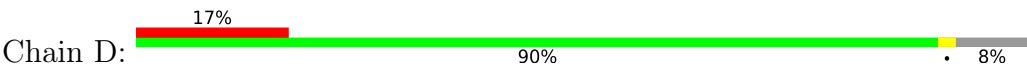
- Molecule 1: Oxidoreductase, zinc-binding dehydrogenase family



- Molecule 1: Oxidoreductase, zinc-binding dehydrogenase family



- Molecule 1: Oxidoreductase, zinc-binding dehydrogenase family



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.29Å 77.43Å 309.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.94 – 2.55 42.94 – 2.55	Depositor EDS
% Data completeness (in resolution range)	100.0 (42.94-2.55) 99.3 (42.94-2.55)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.234 , 0.269 (Not available) , 0.254	Depositor DCC
R_{free} test set	3088 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	53.7	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 65.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.035 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	19587	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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: 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	0.11	0/2633	0.26	0/3582
1	B	0.11	0/2374	0.26	0/3230
1	C	0.11	0/2596	0.26	0/3537
1	D	0.11	0/2390	0.25	0/3248
All	All	0.11	0/9993	0.26	0/13597

There are no bond length outliers.

There are no bond angle outliers.

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	2587	2529	2543	10	1
1	B	2343	2298	2297	10	0
1	C	2553	2480	2481	8	2
1	D	2358	2324	2323	3	0
2	A	2	0	0	0	1
2	B	2	0	0	0	0
2	C	2	0	0	0	2
2	D	2	0	0	0	0
3	A	42	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	18	0	0	0	0
3	C	35	0	0	1	0
3	D	12	0	0	0	0
All	All	9956	9631	9644	30	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 2.

All (30) 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:89:ARG:NH2	1:B:106:CYS:O	2.10	0.84
1:C:75:THR:O	3:C:501:HOH:O	2.01	0.79
1:A:10:GLY:O	1:A:47:ARG:NH1	2.16	0.78
1:C:77:GLY:O	1:C:134:ARG:NH1	2.21	0.73
1:C:24:ARG:NH2	1:C:70:GLU:OE2	2.27	0.67
1:A:77:GLY:O	1:A:134:ARG:NH2	2.29	0.64
1:A:47:ARG:NH2	1:A:330:GLU:OE2	2.33	0.61
1:B:123:GLN:NE2	1:B:343:PRO:O	2.35	0.60
1:A:24:ARG:NH2	1:A:70:GLU:OE2	2.41	0.53
1:B:92:CYS:SG	1:B:108:THR:OG1	2.67	0.52
1:B:97:LEU:O	1:B:102:ARG:N	2.44	0.50
1:B:111:ARG:HD2	1:B:117:LEU:HD21	1.95	0.49
1:D:89:ARG:NH2	1:D:106:CYS:O	2.46	0.48
1:B:116:GLU:OE1	1:B:116:GLU:N	2.42	0.48
1:B:193:ARG:NH2	1:D:213:ASP:OD1	2.42	0.47
1:A:1:MET:HE2	1:A:122:ALA:HB1	1.96	0.47
1:C:-2:GLN:OE1	1:C:24:ARG:NH1	2.49	0.46
1:C:64:ALA:CB	1:C:125:MET:HE1	2.44	0.46
1:C:1:MET:CE	1:C:122:ALA:HB1	2.45	0.46
1:A:228:ARG:O	1:A:233:GLY:N	2.45	0.46
1:B:162:GLY:O	1:B:260:ARG:NH1	2.50	0.45
1:C:92:CYS:SG	1:C:108:THR:OG1	2.76	0.43
1:A:218:SER:OG	1:A:222:ASP:OD1	2.30	0.42
1:D:1:MET:SD	1:D:122:ALA:HB1	2.59	0.42
1:A:-2:GLN:OE1	1:A:24:ARG:NH1	2.46	0.42
1:A:92:CYS:SG	1:A:108:THR:OG1	2.78	0.42
1:B:150:THR:OG1	1:B:151:THR:N	2.53	0.41
1:A:91:SER:OG	1:A:98:CYS:SG	2.79	0.41
1:C:64:ALA:HB3	1:C:125:MET:HE1	2.03	0.40
1:B:117:LEU:N	1:B:117:LEU:HD23	2.36	0.40

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:C:62:HIS:HE2	2:C:402:ZN:ZN[3_545]	1.43	0.17
1:C:149:CYS:HG	2:C:402:ZN:ZN[3_545]	1.54	0.06
1:A:-9:HIS:HE2	2:A:402:ZN:ZN[4_556]	1.56	0.04

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	357/362 (99%)	340 (95%)	17 (5%)	0	100	100
1	B	329/362 (91%)	312 (95%)	17 (5%)	0	100	100
1	C	356/362 (98%)	339 (95%)	16 (4%)	1 (0%)	36	45
1	D	327/362 (90%)	309 (94%)	17 (5%)	1 (0%)	36	45
All	All	1369/1448 (94%)	1300 (95%)	67 (5%)	2 (0%)	48	61

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	-1	GLY
1	D	25	PRO

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	
1	A	261/277 (94%)	259 (99%)	2 (1%)	73	84
1	B	232/277 (84%)	229 (99%)	3 (1%)	61	76
1	C	251/277 (91%)	249 (99%)	2 (1%)	73	84
1	D	236/277 (85%)	233 (99%)	3 (1%)	61	76
All	All	980/1108 (88%)	970 (99%)	10 (1%)	68	80

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	ARG
1	A	275	ASN
1	C	-13	HIS
1	C	-5	LEU
1	B	17	MET
1	B	117	LEU
1	B	267	LEU
1	D	1	MET
1	D	92	CYS
1	D	94	ASP

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	331	ASN
1	C	103	ASN
1	C	273	GLN
1	B	46	ASN
1	B	232	GLN
1	B	294	GLN
1	B	331	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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 [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	358/362 (98%)	0.52	15 (4%)	40 41	30, 68, 109, 132	1 (0%)
1	B	333/362 (91%)	1.33	87 (26%)	1 1	47, 91, 138, 193	0
1	C	358/362 (98%)	0.66	39 (10%)	10 10	45, 73, 113, 149	0
1	D	333/362 (91%)	1.08	63 (18%)	3 2	44, 88, 144, 174	0
All	All	1382/1448 (95%)	0.89	204 (14%)	5 5	30, 78, 130, 193	1 (0%)

All (204) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	320	ALA	7.2
1	B	315	ASP	6.4
1	B	3	ALA	5.8
1	D	49	GLY	5.6
1	D	15	LEU	5.3
1	A	333	GLU	5.2
1	D	315	ASP	5.1
1	B	6	LEU	4.7
1	B	312	PHE	4.5
1	A	336	GLY	4.4
1	D	335	ALA	4.4
1	D	10	GLY	4.4
1	B	320	ALA	4.4
1	B	334	VAL	4.3
1	A	332	ARG	4.3
1	D	7	ARG	4.2
1	C	205	GLN	4.2
1	C	344	GLU	4.1
1	C	308	GLY	4.1
1	D	332	ARG	4.0
1	C	176	SER	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	267	LEU	4.0
1	D	91	SER	3.9
1	B	1	MET	3.9
1	B	316	ALA	3.9
1	B	333	GLU	3.9
1	D	9	PHE	3.9
1	B	76	PRO	3.9
1	D	337	ARG	3.8
1	B	94	ASP	3.8
1	C	-13	HIS	3.8
1	B	121	TYR	3.8
1	C	337	ARG	3.7
1	B	17	MET	3.7
1	A	205	GLN	3.6
1	B	327	PHE	3.6
1	D	4	VAL	3.5
1	D	118	PRO	3.5
1	B	337	ARG	3.5
1	A	201	GLU	3.5
1	B	20	MET	3.5
1	A	315	ASP	3.5
1	B	339	VAL	3.5
1	B	329	LEU	3.5
1	B	0	HIS	3.5
1	B	93	GLY	3.4
1	D	108	THR	3.4
1	A	-13	HIS	3.4
1	B	268	GLU	3.4
1	D	222	ASP	3.4
1	D	47	ARG	3.4
1	D	104	SER	3.4
1	B	340	LEU	3.3
1	B	332	ARG	3.3
1	B	37	CYS	3.3
1	B	11	GLU	3.3
1	A	337	ARG	3.3
1	B	4	VAL	3.3
1	B	341	VAL	3.3
1	C	216	ILE	3.2
1	B	12	ALA	3.2
1	A	219	ARG	3.2
1	D	266	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	73	PRO	3.2
1	D	103	ASN	3.2
1	B	34	VAL	3.2
1	B	71	VAL	3.2
1	B	77	GLY	3.1
1	C	306	THR	3.1
1	B	24	ARG	3.1
1	D	333	GLU	3.1
1	C	223	PHE	3.1
1	C	200	SER	3.1
1	D	14	ASN	3.1
1	C	232	GLN	3.1
1	C	-12	HIS	3.1
1	B	50	ASN	3.1
1	C	335	ALA	3.0
1	B	96	ALA	3.0
1	C	312	PHE	3.0
1	B	7	ARG	3.0
1	B	66	GLY	3.0
1	C	217	VAL	3.0
1	B	124	PHE	3.0
1	B	313	VAL	3.0
1	D	117	LEU	3.0
1	D	2	LYS	3.0
1	B	343	PRO	2.9
1	B	247	THR	2.9
1	D	99	ARG	2.9
1	B	97	LEU	2.9
1	D	97	LEU	2.9
1	D	43	ASP	2.9
1	B	126	VAL	2.9
1	D	314	THR	2.9
1	B	325	ALA	2.9
1	D	8	SER	2.8
1	C	221	LEU	2.8
1	B	41	TYR	2.8
1	B	129	VAL	2.8
1	C	203	LYS	2.8
1	B	45	ILE	2.8
1	D	269	SER	2.8
1	B	32	LEU	2.7
1	C	75	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	19	THR	2.7
1	B	8	SER	2.7
1	B	271	MET	2.7
1	B	219	ARG	2.7
1	B	9	PHE	2.6
1	B	143	GLU	2.6
1	B	2	LYS	2.6
1	D	21	PRO	2.6
1	C	213	ASP	2.6
1	B	63	GLU	2.6
1	B	88	GLN	2.6
1	C	229	LYS	2.6
1	B	28	GLY	2.6
1	D	76	PRO	2.6
1	B	322	ALA	2.6
1	D	127	ALA	2.6
1	D	287	LEU	2.5
1	C	202	ALA	2.5
1	C	77	GLY	2.5
1	D	109	ASP	2.5
1	C	273	GLN	2.5
1	C	169	THR	2.5
1	B	106	CYS	2.5
1	C	211	GLY	2.5
1	D	5	VAL	2.5
1	D	325	ALA	2.5
1	A	191	GLY	2.5
1	D	120	GLY	2.5
1	B	27	ARG	2.5
1	D	46	ASN	2.5
1	B	189	LEU	2.4
1	D	48	ARG	2.4
1	B	39	VAL	2.4
1	B	145	ALA	2.4
1	C	336	GLY	2.4
1	B	120	GLY	2.4
1	B	211	GLY	2.4
1	C	315	ASP	2.4
1	B	69	ILE	2.4
1	D	6	LEU	2.4
1	C	219	ARG	2.4
1	C	228	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	99	ARG	2.4
1	B	22	MET	2.4
1	C	234	ALA	2.4
1	B	314	THR	2.4
1	B	25	PRO	2.4
1	B	130	GLY	2.4
1	C	41	TYR	2.4
1	A	334	VAL	2.4
1	B	107	LYS	2.4
1	D	202	ALA	2.4
1	B	49	GLY	2.3
1	B	91	SER	2.3
1	D	338	LEU	2.3
1	A	335	ALA	2.3
1	D	246	ALA	2.3
1	D	201	GLU	2.3
1	D	323	ALA	2.3
1	D	327	PHE	2.3
1	C	146	THR	2.2
1	D	130	GLY	2.2
1	A	202	ALA	2.2
1	B	92	CYS	2.2
1	B	108	THR	2.2
1	B	265	GLY	2.2
1	C	201	GLU	2.2
1	D	268	GLU	2.2
1	C	324	LYS	2.2
1	D	90	MET	2.2
1	D	329	LEU	2.2
1	C	310	ARG	2.2
1	B	21	PRO	2.2
1	A	235	GLY	2.2
1	D	-1	GLY	2.2
1	C	338	LEU	2.1
1	B	338	LEU	2.1
1	D	60	LEU	2.1
1	D	87	LEU	2.1
1	D	92	CYS	2.1
1	B	175	ALA	2.1
1	D	3	ALA	2.1
1	C	233	GLY	2.1
1	D	143	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	294	GLN	2.1
1	B	64	ALA	2.1
1	B	122	ALA	2.1
1	D	13	GLY	2.1
1	D	114	GLY	2.1
1	A	190	ASP	2.1
1	D	0	HIS	2.1
1	C	188	ARG	2.0
1	C	281	VAL	2.0
1	B	68	VAL	2.0
1	B	83	ARG	2.0
1	D	27	ARG	2.0
1	D	113	PHE	2.0
1	B	10	GLY	2.0
1	D	62	HIS	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	ZN	B	401	1/1	0.87	0.18	108,108,108,108	0
2	ZN	D	401	1/1	0.93	0.12	117,117,117,117	0
2	ZN	C	402	1/1	0.94	0.16	89,89,89,89	0
2	ZN	B	402	1/1	0.95	0.06	107,107,107,107	0
2	ZN	D	402	1/1	0.95	0.07	107,107,107,107	0
2	ZN	A	402	1/1	0.96	0.11	70,70,70,70	0
2	ZN	A	401	1/1	0.97	0.11	67,67,67,67	0
2	ZN	C	401	1/1	0.99	0.09	61,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.