



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 9, 2026 – 04:10 PM JST

PDB ID : 9WVJ / pdb_00009wvj
Title : crystal structure of an anti-phage protein complex
Authors : Wang, H.; Feng, Y.; Liu, X.
Deposited on : 2025-09-20
Resolution : 2.09 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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

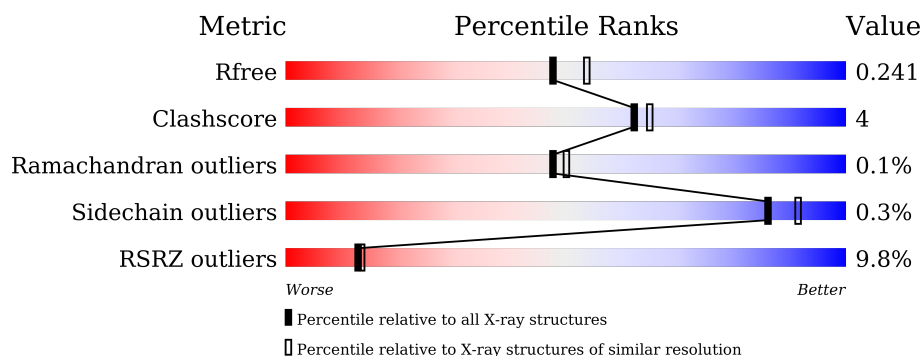
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	471	6% 87% 8% .
1	B	471	6% 90% 6% .
2	C	274	15% 83% 8% 8%
2	D	274	15% 81% 9% 9% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11404 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 7-cyano-7-deazaguanine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	454	Total	C	N	O	S	0	0	0
			3533	2213	638	666	16			
1	A	452	Total	C	N	O	S	0	0	0
			3522	2207	636	663	16			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	MET	-	initiating methionine	UNP A0A4Q0WMG2
B	-12	GLY	-	expression tag	UNP A0A4Q0WMG2
B	-11	SER	-	expression tag	UNP A0A4Q0WMG2
B	-10	SER	-	expression tag	UNP A0A4Q0WMG2
B	-9	HIS	-	expression tag	UNP A0A4Q0WMG2
B	-8	HIS	-	expression tag	UNP A0A4Q0WMG2
B	-7	HIS	-	expression tag	UNP A0A4Q0WMG2
B	-6	HIS	-	expression tag	UNP A0A4Q0WMG2
B	-5	HIS	-	expression tag	UNP A0A4Q0WMG2
B	-4	HIS	-	expression tag	UNP A0A4Q0WMG2
B	-3	SER	-	expression tag	UNP A0A4Q0WMG2
B	-2	GLN	-	expression tag	UNP A0A4Q0WMG2
B	-1	ASP	-	expression tag	UNP A0A4Q0WMG2
B	0	PRO	-	expression tag	UNP A0A4Q0WMG2
A	-13	MET	-	initiating methionine	UNP A0A4Q0WMG2
A	-12	GLY	-	expression tag	UNP A0A4Q0WMG2
A	-11	SER	-	expression tag	UNP A0A4Q0WMG2
A	-10	SER	-	expression tag	UNP A0A4Q0WMG2
A	-9	HIS	-	expression tag	UNP A0A4Q0WMG2
A	-8	HIS	-	expression tag	UNP A0A4Q0WMG2
A	-7	HIS	-	expression tag	UNP A0A4Q0WMG2
A	-6	HIS	-	expression tag	UNP A0A4Q0WMG2
A	-5	HIS	-	expression tag	UNP A0A4Q0WMG2
A	-4	HIS	-	expression tag	UNP A0A4Q0WMG2
A	-3	SER	-	expression tag	UNP A0A4Q0WMG2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLN	-	expression tag	UNP A0A4Q0WMG2
A	-1	ASP	-	expression tag	UNP A0A4Q0WMG2
A	0	PRO	-	expression tag	UNP A0A4Q0WMG2

- Molecule 2 is a protein called DUF5610 domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	250	Total	C	N	O	S	0	0	0
			1860	1153	328	373	6			
2	C	251	Total	C	N	O	S	0	0	0
			1869	1157	331	376	5			

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

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

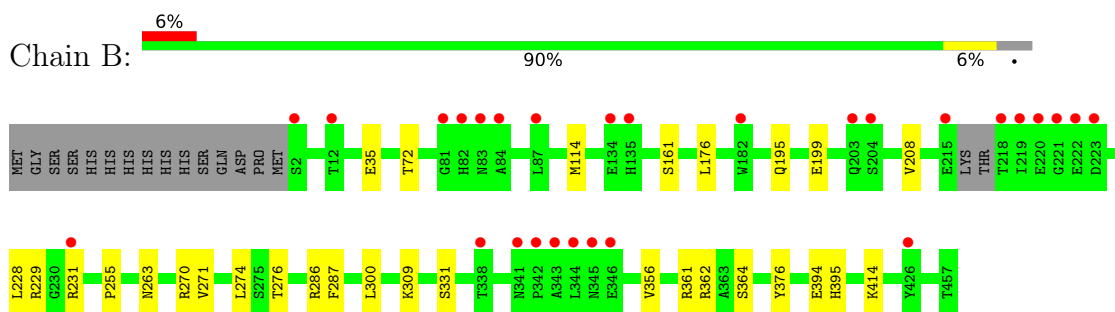
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	251	Total	O	0	0
			251	251		
4	D	108	Total	O	0	0
			108	108		
4	A	176	Total	O	0	0
			176	176		
4	C	83	Total	O	0	0
			83	83		

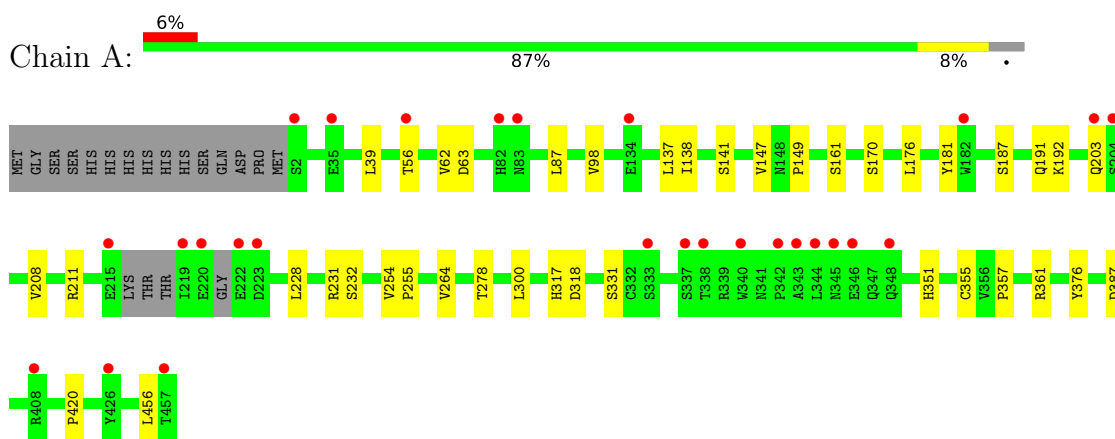
3 Residue-property plots [i](#)

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.

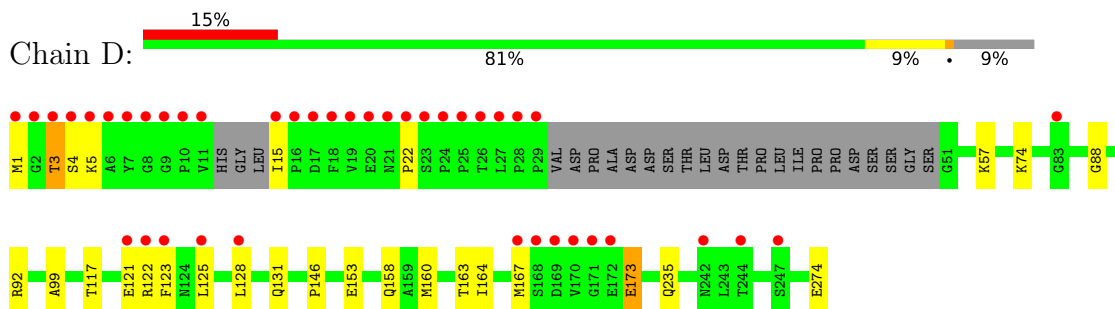
- Molecule 1: 7-cyano-7-deazaguanine synthase



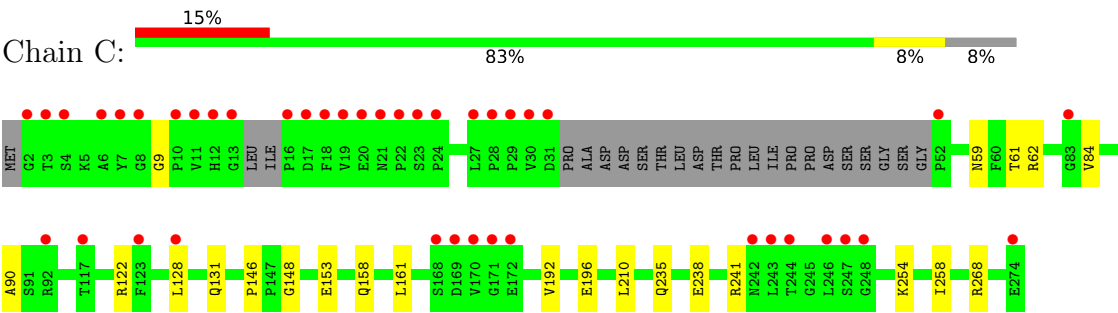
- Molecule 1: 7-cyano-7-deazaguanine synthase



- Molecule 2: DUF5610 domain-containing protein



- Molecule 2: DUF5610 domain-containing protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.39Å 116.18Å 120.32Å 90.00° 103.01° 90.00°	Depositor
Resolution (Å)	58.62 – 2.09 58.62 – 2.09	Depositor EDS
% Data completeness (in resolution range)	99.5 (58.62-2.09) 99.7 (58.62-2.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.08Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.206 , 0.241 0.206 , 0.241	Depositor DCC
R_{free} test set	4529 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	20.7	Xtriage
Anisotropy	0.790	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.022 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11404	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.12% 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	0.17	0/3598	0.37	0/4880
1	B	0.18	0/3610	0.39	0/4898
2	C	0.22	0/1896	0.43	0/2558
2	D	0.23	0/1886	0.49	0/2545
All	All	0.19	0/10990	0.41	0/14881

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	3522	0	3465	26	0
1	B	3533	0	3476	28	0
2	C	1869	0	1836	15	0
2	D	1860	0	1837	29	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	176	0	0	1	0
4	B	251	0	0	3	0
4	C	83	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	108	0	0	1	0
All	All	11404	0	10614	87	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 4.

All (87) 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)
2:D:74:LYS:HG2	2:D:274:GLU:HG2	1.62	0.80
1:B:395:HIS:CE1	2:D:4:SER:HA	2.21	0.75
1:B:263:ASN:HD21	2:D:3:THR:HB	1.53	0.71
2:D:125:LEU:HB3	2:D:128:LEU:HD23	1.73	0.69
1:A:228:LEU:O	1:A:231:ARG:HD3	1.92	0.69
1:B:35:GLU:H	1:B:35:GLU:CD	2.00	0.69
1:B:276:THR:H	2:D:3:THR:HA	1.59	0.68
1:B:195:GLN:O	1:B:199:GLU:HG3	1.93	0.67
1:A:147:VAL:HG23	1:A:149:PRO:HD3	1.75	0.67
1:A:62:VAL:HG11	1:A:138:ILE:HD11	1.79	0.64
1:B:228:LEU:HD13	1:B:231:ARG:HD3	1.78	0.64
1:A:56:THR:HG21	1:A:141:SER:H	1.63	0.63
1:A:264:VAL:HA	1:A:420:PRO:HG3	1.81	0.61
2:D:88:GLY:O	2:D:92:ARG:HG2	2.02	0.60
2:C:238:GLU:O	2:C:241:ARG:HD2	2.02	0.60
1:A:203:GLN:H	1:A:203:GLN:CD	2.09	0.60
2:D:3:THR:HG23	2:D:4:SER:H	1.68	0.59
1:B:395:HIS:HE1	2:D:4:SER:HA	1.65	0.59
2:D:125:LEU:HD12	2:D:128:LEU:HD21	1.86	0.58
1:A:176:LEU:HD11	1:A:208:VAL:HG23	1.86	0.58
1:A:161:SER:HA	1:A:255:PRO:HG2	1.87	0.57
1:B:276:THR:HB	2:D:1:MET:HE3	1.87	0.56
1:A:56:THR:CG2	1:A:141:SER:H	2.18	0.55
2:C:59:ASN:ND2	2:C:62:ARG:HH21	2.05	0.55
1:B:331:SER:HB2	1:B:361:ARG:HD2	1.88	0.55
1:B:287:PHE:CD1	1:B:300:LEU:HD11	2.42	0.55
1:B:395:HIS:HD2	4:B:603:HOH:O	1.89	0.55
2:C:128:LEU:HA	2:C:131:GLN:HG3	1.89	0.55
4:B:767:HOH:O	2:D:4:SER:HB2	2.08	0.53
1:B:270:ARG:NH2	1:B:394:GLU:HG3	2.23	0.53
1:B:274:LEU:HD12	2:D:5:LYS:HG3	1.90	0.53
1:B:270:ARG:HH22	1:B:394:GLU:HG3	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:158:GLN:HG2	4:D:387:HOH:O	2.09	0.52
2:D:22:PRO:HG2	2:D:99:ALA:HB2	1.93	0.51
1:A:232:SER:HB2	1:A:278:THR:HG23	1.93	0.51
1:B:309:LYS:HB3	1:B:364:SER:HB3	1.93	0.51
2:C:254:LYS:O	2:C:258:ILE:HG12	2.11	0.50
1:B:263:ASN:ND2	2:D:3:THR:HB	2.25	0.50
2:D:146:PRO:HG2	2:D:153:GLU:HG2	1.94	0.50
1:A:181:TYR:CD2	1:A:187:SER:HB2	2.47	0.49
1:B:161:SER:HA	1:B:255:PRO:HG2	1.95	0.49
2:D:117:THR:O	2:D:121:GLU:HG2	2.12	0.48
1:A:39:LEU:HD21	1:A:137:LEU:HD11	1.95	0.48
2:D:121:GLU:O	2:D:122:ARG:HB3	2.14	0.48
1:A:181:TYR:O	1:A:211:ARG:HA	2.14	0.48
1:B:356:VAL:HG21	2:D:3:THR:HG23	1.95	0.48
2:D:121:GLU:C	2:D:123:PHE:H	2.21	0.47
2:C:148:GLY:HA2	2:C:153:GLU:HB3	1.96	0.47
1:A:170:SER:HB3	1:A:317:HIS:HD2	1.80	0.47
1:A:170:SER:HB3	1:A:317:HIS:CD2	2.50	0.46
2:C:122:ARG:HE	2:C:122:ARG:HB2	1.57	0.46
2:D:164:ILE:O	2:D:167:MET:HG2	2.15	0.46
1:B:114:MET:HG3	1:B:286:ARG:CB	2.46	0.46
1:B:176:LEU:HD11	1:B:208:VAL:HG23	1.99	0.45
2:D:57:LYS:HE3	2:D:57:LYS:HB2	1.72	0.45
1:A:181:TYR:CZ	1:A:211:ARG:HB2	2.52	0.45
2:C:235:GLN:H	2:C:235:GLN:CD	2.23	0.45
2:D:235:GLN:CD	2:D:235:GLN:H	2.25	0.44
1:B:414:LYS:HE2	4:B:834:HOH:O	2.17	0.44
1:B:362:ARG:HG2	1:B:376:TYR:CZ	2.52	0.44
1:A:387:ASP:CG	1:A:456:LEU:HB2	2.43	0.44
1:B:356:VAL:HG21	2:D:3:THR:CG2	2.48	0.44
2:C:192:VAL:HG12	2:C:196:GLU:OE2	2.18	0.44
1:A:63:ASP:HB3	1:A:98:VAL:HG12	2.00	0.43
1:A:203:GLN:H	1:A:203:GLN:NE2	2.16	0.43
1:A:170:SER:OG	1:A:318:ASP:HB2	2.18	0.43
1:B:229:ARG:HA	2:D:1:MET:HE1	2.01	0.43
1:B:271:VAL:O	1:B:274:LEU:HD23	2.18	0.43
1:B:72:THR:HG23	1:B:231:ARG:NH1	2.34	0.43
1:A:254:VAL:HG21	1:A:300:LEU:HD23	2.00	0.43
1:B:114:MET:HG3	1:B:286:ARG:HB2	2.00	0.42
1:A:87:LEU:HD12	2:C:9:GLY:O	2.20	0.42
2:C:84:VAL:HG21	2:C:90:ALA:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:268:ARG:HD2	2:C:268:ARG:HA	1.91	0.42
1:A:351:HIS:HB2	1:A:376:TYR:CD1	2.55	0.42
2:D:128:LEU:HA	2:D:131:GLN:HG3	2.00	0.42
1:A:355:CYS:SG	1:A:357:PRO:HD2	2.60	0.42
2:D:160:MET:O	2:D:163:THR:HG22	2.20	0.41
2:D:125:LEU:HD12	2:D:128:LEU:CD2	2.50	0.41
2:D:173:GLU:H	2:D:173:GLU:HG3	1.67	0.41
1:A:191:GLN:HG3	1:A:192:LYS:N	2.35	0.41
2:C:59:ASN:HD21	2:C:62:ARG:HH21	1.67	0.41
2:C:61:THR:HG23	2:C:210:LEU:HD11	2.02	0.41
1:B:72:THR:HG23	1:B:231:ARG:HH12	1.86	0.40
2:C:146:PRO:HG2	2:C:153:GLU:HG2	2.01	0.40
4:A:769:HOH:O	2:C:158:GLN:HB2	2.20	0.40
1:A:331:SER:HB2	1:A:361:ARG:HD2	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	
1	A	446/471 (95%)	439 (98%)	7 (2%)	0	100	100
1	B	450/471 (96%)	444 (99%)	6 (1%)	0	100	100
2	C	245/274 (89%)	238 (97%)	7 (3%)	0	100	100
2	D	244/274 (89%)	236 (97%)	7 (3%)	1 (0%)	30	28
All	All	1385/1490 (93%)	1357 (98%)	27 (2%)	1 (0%)	48	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	3	THR

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	385/402 (96%)	385 (100%)	0	100	100
1	B	386/402 (96%)	386 (100%)	0	100	100
2	C	196/216 (91%)	195 (100%)	1 (0%)	81	88
2	D	195/216 (90%)	193 (99%)	2 (1%)	68	76
All	All	1162/1236 (94%)	1159 (100%)	3 (0%)	86	91

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	15	ILE
2	D	173	GLU
2	C	161	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	36	ASN
1	B	107	GLN
1	B	126	HIS
1	B	191	GLN
1	B	263	ASN
1	B	293	ASN
1	B	395	HIS
1	B	400	GLN
2	D	112	GLN
1	A	36	ASN
1	A	82	HIS
1	A	126	HIS
1	A	203	GLN
1	A	253	ASN
1	A	299	HIS
1	A	324	GLN

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Mol	Chain	Res	Type
1	A	400	GLN
2	C	59	ASN
2	C	127	ASN
2	C	131	GLN
2	C	158	GLN
2	C	206	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 ⓘ

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

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	452/471 (95%)	0.51	27 (5%)	27	29	18, 30, 51, 97	1 (0%)
1	B	454/471 (96%)	0.23	28 (6%)	26	28	11, 21, 47, 86	0
2	C	251/274 (91%)	0.80	42 (16%)	4	4	16, 32, 71, 97	0
2	D	250/274 (91%)	0.82	41 (16%)	4	4	12, 27, 78, 100	1 (0%)
All	All	1407/1490 (94%)	0.53	138 (9%)	13	13	11, 27, 60, 100	2 (0%)

All (138) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	3	THR	11.6
2	D	2	GLY	8.6
2	D	5	LYS	8.0
2	D	4	SER	7.9
2	D	22	PRO	7.1
2	D	1	MET	6.7
2	D	19	VAL	6.2
2	C	30	VAL	6.2
2	D	6	ALA	6.2
2	D	18	PHE	6.2
2	D	10	PRO	5.7
2	C	2	GLY	5.5
2	D	11	VAL	5.5
2	D	7	TYR	4.8
2	D	168	SER	4.7
2	D	23	SER	4.6
1	B	219	ILE	4.5
2	C	13	GLY	4.5
2	D	16	PRO	4.5
2	C	19	VAL	4.4
1	A	219	ILE	4.4

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Mol	Chain	Res	Type	RSRZ
2	C	10	PRO	4.4
1	A	220	GLU	4.4
1	B	221	GLY	4.3
2	C	18	PHE	4.2
1	B	220	GLU	4.2
1	B	342	PRO	4.2
2	C	31	ASP	4.2
1	A	222	GLU	4.1
2	D	15	ILE	4.1
1	A	182	TRP	4.0
2	C	12	HIS	4.0
2	D	170	VAL	4.0
2	D	24	PRO	4.0
1	B	82	HIS	3.9
1	B	345	ASN	3.9
2	D	122	ARG	3.9
1	B	215	GLU	3.8
1	B	344	LEU	3.8
2	C	29	PRO	3.7
2	C	247	SER	3.7
2	C	22	PRO	3.7
2	C	172	GLU	3.7
2	C	3	THR	3.7
2	D	172	GLU	3.6
1	B	343	ALA	3.6
1	B	218	THR	3.5
2	D	9	GLY	3.5
2	C	11	VAL	3.5
2	C	6	ALA	3.4
2	D	247	SER	3.4
2	D	121	GLU	3.4
2	D	29	PRO	3.4
1	A	340	TRP	3.3
1	A	343	ALA	3.3
2	C	28	PRO	3.3
2	D	26	THR	3.2
2	C	170	VAL	3.2
1	A	223	ASP	3.2
1	A	333	SER	3.2
1	B	222	GLU	3.2
1	A	344	LEU	3.2
2	C	7	TYR	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	82	HIS	3.1
2	C	16	PRO	3.1
2	D	244	THR	3.1
2	C	17	ASP	3.1
2	C	169	ASP	3.1
2	C	20	GLU	3.0
1	A	203	GLN	3.0
2	D	8	GLY	2.9
1	A	204	SER	2.9
2	D	27	LEU	2.9
2	D	28	PRO	2.9
2	C	8	GLY	2.9
1	B	203	GLN	2.8
2	C	21	ASN	2.8
1	B	182	TRP	2.8
2	C	274	GLU	2.8
1	B	341	ASN	2.8
1	A	337	SER	2.8
2	D	169	ASP	2.7
1	B	338	THR	2.7
2	C	4	SER	2.6
2	D	83	GLY	2.6
1	A	346	GLU	2.6
2	D	20	GLU	2.6
1	B	87	LEU	2.6
1	A	345	ASN	2.6
2	C	52	PRO	2.6
1	A	342	PRO	2.5
2	D	123	PHE	2.5
1	B	83	ASN	2.5
2	D	242	ASN	2.5
2	D	125	LEU	2.5
1	A	83	ASN	2.5
1	B	223	ASP	2.4
1	A	215	GLU	2.4
2	C	246	LEU	2.4
1	B	204	SER	2.4
2	D	167	MET	2.4
2	C	24	PRO	2.4
1	A	134	GLU	2.4
2	C	27	LEU	2.4
1	B	2	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	56	THR	2.4
2	C	123	PHE	2.3
1	A	338	THR	2.3
1	B	81	GLY	2.3
2	C	168	SER	2.3
2	C	242	ASN	2.3
1	B	84	ALA	2.3
2	D	171	GLY	2.3
2	C	171	GLY	2.3
1	B	135	HIS	2.2
1	A	2	SER	2.2
2	C	244	THR	2.2
2	D	21	ASN	2.2
2	C	243	LEU	2.2
1	B	346	GLU	2.2
1	A	457	THR	2.2
2	D	25	PRO	2.2
2	C	117	THR	2.2
2	C	248	GLY	2.2
1	A	348	GLN	2.2
1	B	231	ARG	2.2
2	D	17	ASP	2.2
1	B	426	TYR	2.1
2	D	128	LEU	2.1
2	C	128	LEU	2.1
2	C	83	GLY	2.1
2	C	23	SER	2.1
1	A	35	GLU	2.0
1	A	408	ARG	2.0
2	C	92	ARG	2.0
1	A	426	TYR	2.0
1	B	134	GLU	2.0
1	B	12	THR	2.0

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

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

6.3 Carbohydrates ⓘ

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
3	ZN	B	500	1/1	0.98	0.04	19,19,19,19	0
3	ZN	A	500	1/1	0.99	0.02	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.