



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:47 AM UTC

PDB ID : 7ZMT / pdb_00007zmt
Title : Crystal structure of human RECQL5 helicase APO form in complex with engineered nanobody (Gluebody) G5-006
Authors : Ye, M.; Makola, M.; Newman, J.A.; Fairhead, M.; MacLean, E.; Krojer, T.; Aitkenhead, H.; Bountra, C.; Gileadi, O.; von Delft, F.
Deposited on : 2022-04-19
Resolution : 2.30 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

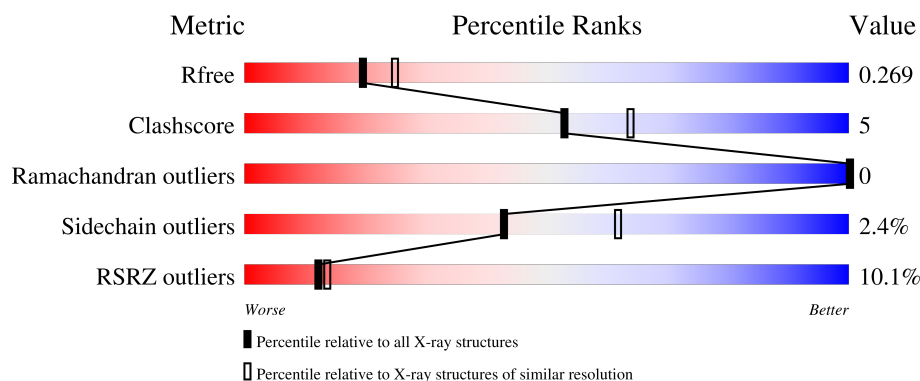
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	445	
1	B	445	
2	C	127	
2	K	127	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8893 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 ATP-dependent DNA helicase Q5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	443	Total	C	N	O	S	0	3	0
			3454	2180	621	631	22			
1	B	441	Total	C	N	O	S	0	3	0
			3446	2175	620	628	23			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	SER	-	expression tag	UNP O94762
A	10	MET	-	expression tag	UNP O94762
B	9	SER	-	expression tag	UNP O94762
B	10	MET	-	expression tag	UNP O94762

- Molecule 2 is a protein called Gluebody G5-006.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	124	Total	C	N	O	S	0	2	0
			954	591	164	193	6			
2	C	124	Total	C	N	O	S	0	2	0
			954	591	164	193	6			

- 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		

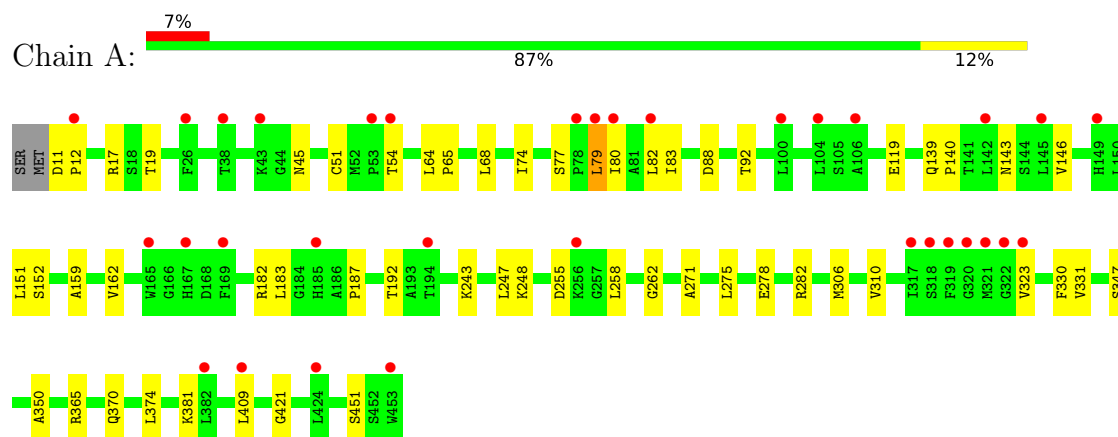
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total 32	O 32	0	0
4	B	31	Total 31	O 31	0	0
4	K	12	Total 12	O 12	0	0
4	C	8	Total 8	O 8	0	0

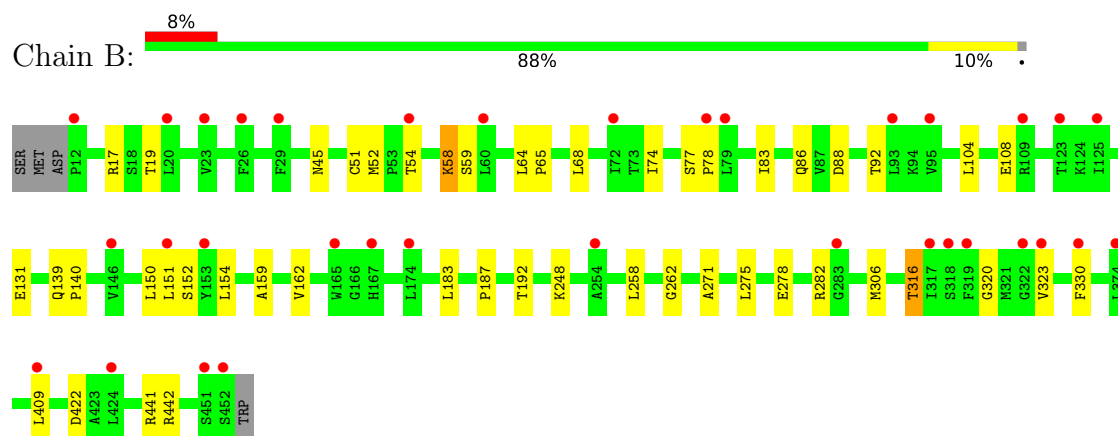
3 Residue-property plots [i](#)

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.

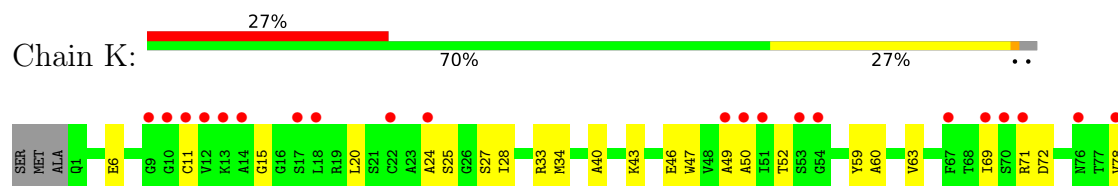
- Molecule 1: ATP-dependent DNA helicase Q5

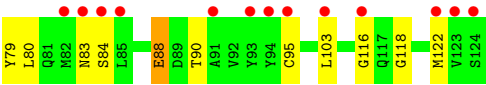


- Molecule 1: ATP-dependent DNA helicase Q5

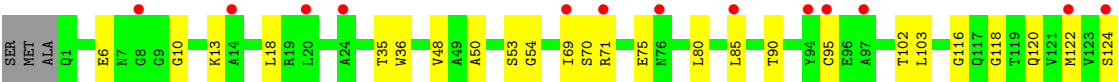
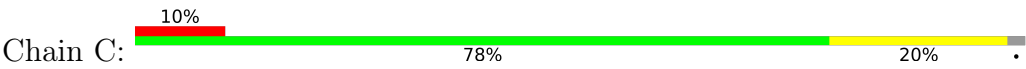


- Molecule 2: Gluebody G5-006





● Molecule 2: Gluebody G5-006



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	82.36Å 90.67Å 217.85Å 90.00° 90.75° 90.00°	Depositor
Resolution (Å)	56.74 – 2.30 56.74 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (56.74-2.30) 99.7 (56.74-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.264 , 0.291 0.257 , 0.269	Depositor DCC
R_{free} test set	2761 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	58.7	Xtriage
Anisotropy	0.183	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 36.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.190 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8893	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.07% 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: 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	1.10	1/3527 (0.0%)	1.53	1/4768 (0.0%)
1	B	1.11	3/3516 (0.1%)	1.54	10/4751 (0.2%)
2	C	1.13	0/979	1.39	2/1328 (0.2%)
2	K	1.27	4/979 (0.4%)	1.36	11/1328 (0.8%)
All	All	1.12	8/9001 (0.1%)	1.50	24/12175 (0.2%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	K	83[A]	ASN	C-O	11.16	1.38	1.23
2	K	83[B]	ASN	C-O	11.16	1.38	1.23
2	K	71[A]	ARG	C-O	10.55	1.36	1.24
2	K	71[B]	ARG	C-O	10.55	1.36	1.24
1	B	78	PRO	C-N	6.61	1.42	1.33
1	B	441[A]	ARG	C-O	5.29	1.30	1.24
1	B	441[B]	ARG	C-O	5.29	1.30	1.24
1	A	310	VAL	N-CA	5.25	1.50	1.46

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	71[A]	ARG	CA-C-O	10.15	131.32	120.36
2	K	71[B]	ARG	CA-C-O	10.15	131.32	120.36
2	K	71[A]	ARG	N-CA-C	8.20	122.27	108.90
2	K	71[B]	ARG	N-CA-C	8.20	122.27	108.90
1	B	78	PRO	O-C-N	7.17	131.19	122.23
2	K	71[A]	ARG	O-C-N	-6.84	115.16	123.30
2	K	71[B]	ARG	O-C-N	-6.84	115.16	123.30
2	K	83[A]	ASN	CA-C-O	6.13	128.64	121.46
2	K	83[B]	ASN	CA-C-O	6.13	128.64	121.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	58	LYS	CB-CA-C	-6.09	100.33	110.68
2	C	118	GLY	CA-C-O	-6.02	118.08	122.23
2	K	83[A]	ASN	N-CA-C	5.86	119.11	109.85
2	K	83[B]	ASN	N-CA-C	5.86	119.11	109.85
1	A	54	THR	CB-CA-C	5.62	118.27	109.84
1	B	54	THR	CB-CA-C	5.51	118.49	109.90
2	K	118	GLY	CA-C-O	-5.50	118.43	122.23
2	C	70	SER	CA-C-O	-5.49	115.56	121.38
1	B	316	THR	CA-C-N	5.38	128.12	120.53
1	B	316	THR	C-N-CA	5.38	128.12	120.53
1	B	108	GLU	O-C-N	5.28	127.72	122.12
1	B	17	ARG	CA-C-N	5.16	127.96	120.79
1	B	17	ARG	C-N-CA	5.16	127.96	120.79
1	B	131	GLU	CA-C-N	5.14	127.17	120.28
1	B	131	GLU	C-N-CA	5.14	127.17	120.28

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	3454	0	3469	34	0
1	B	3446	0	3467	21	0
2	C	954	0	900	17	0
2	K	954	0	900	18	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	32	0	0	5	0
4	B	31	0	0	0	0
4	C	8	0	0	2	0
4	K	12	0	0	1	0
All	All	8893	0	8736	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 5.

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)
1:B:306:MET:HE1	1:B:323:VAL:HG11	1.43	0.98
2:C:75:GLU:HA	4:C:205:HOH:O	1.74	0.86
1:A:243:LYS:HD3	1:A:282:ARG:HB3	1.76	0.67
1:A:79:LEU:HG	1:A:82:LEU:HD12	1.77	0.66
2:C:75:GLU:HG2	4:C:205:HOH:O	1.95	0.66
2:C:54:GLY:H	2:C:71[B]:ARG:HH22	1.47	0.63
2:C:50:ALA:C	2:C:69:ILE:HD11	2.24	0.62
1:A:17:ARG:HD2	4:A:603:HOH:O	1.99	0.61
1:A:64:LEU:C	1:A:64:LEU:HD23	2.25	0.61
2:K:90:THR:HG23	2:K:122:MET:HA	1.82	0.60
1:B:64:LEU:C	1:B:64:LEU:HD23	2.27	0.59
1:A:365:ARG:HH12	1:A:451:SER:HB2	1.68	0.58
1:A:365:ARG:NH1	1:A:451:SER:HB2	2.19	0.57
1:A:306:MET:HE1	1:A:323:VAL:HG13	1.87	0.56
1:A:88:ASP:O	1:A:92:THR:HG23	2.05	0.56
1:B:88:ASP:O	1:B:92:THR:HG23	2.06	0.55
1:B:159:ALA:O	1:B:162:VAL:HG12	2.07	0.54
1:A:331:VAL:HG11	1:A:347:SER:HA	1.88	0.54
2:C:18:LEU:HD12	2:C:85:LEU:CD1	2.38	0.54
1:B:422:ASP:HA	2:K:103:LEU:HD23	1.90	0.53
2:C:90:THR:HG23	2:C:122:MET:HA	1.90	0.53
1:A:159:ALA:O	1:A:162:VAL:HG12	2.09	0.53
2:K:33:ARG:HD2	2:K:50:ALA:HB1	1.92	0.52
2:C:53:SER:HA	2:C:71[B]:ARG:HH12	1.75	0.52
1:B:316:THR:O	1:B:320:GLY:N	2.35	0.51
1:A:11:ASP:HB2	1:A:12:PRO:HD2	1.93	0.50
1:B:139:GLN:N	1:B:140:PRO:HD2	2.27	0.50
1:A:331:VAL:HG23	1:A:350:ALA:HB2	1.94	0.49
1:A:143:ASN:O	1:A:146:VAL:HG22	2.13	0.49
1:A:139:GLN:N	1:A:140:PRO:HD2	2.27	0.49
2:C:13:LYS:HA	2:C:124:SER:OG	2.12	0.48
2:K:24:ALA:HB1	2:K:28:ILE:HD12	1.96	0.48
1:B:306:MET:HE1	1:B:323:VAL:CG1	2.30	0.48
1:A:74:ILE:HG13	1:A:151:LEU:HD11	1.96	0.48
1:A:271:ALA:O	1:A:275:LEU:HB2	2.14	0.47
1:A:45:ASN:OD1	1:A:45:ASN:N	2.45	0.47
2:K:20:LEU:HD12	2:K:80:LEU:HD23	1.97	0.47
2:K:60:ALA:HB3	2:K:63:VAL:HG22	1.97	0.47
1:A:151:LEU:HD21	1:A:183:LEU:HD13	1.97	0.47
1:B:74:ILE:HG13	1:B:151:LEU:HD11	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:278:GLU:OE1	1:B:282:ARG:NH1	2.47	0.47
1:A:19:THR:OG1	1:A:68:LEU:HD21	2.16	0.46
1:B:271:ALA:O	1:B:275:LEU:HB2	2.15	0.46
1:B:52:MET:HB2	1:B:58:LYS:HG2	1.98	0.46
1:B:51:CYS:HA	1:B:192:THR:O	2.16	0.46
2:K:49:ALA:HB1	2:K:69:ILE:HD13	1.96	0.46
1:A:17:ARG:NH1	4:A:603:HOH:O	2.49	0.45
1:A:51:CYS:HA	1:A:192:THR:O	2.15	0.45
1:A:262:GLY:HA3	1:A:330:PHE:CE2	2.52	0.45
1:B:19:THR:OG1	1:B:68:LEU:HD21	2.17	0.45
2:K:72:ASP:HB2	2:K:79:TYR:HE2	1.81	0.45
2:C:36:TRP:CZ2	2:C:95:CYS:SG	3.09	0.45
2:K:6:GLU:OE2	2:K:116:GLY:HA3	2.15	0.45
2:K:34:MET:HG2	2:K:78:VAL:HG11	1.98	0.45
2:K:88:GLU:H	2:K:88:GLU:HG3	1.52	0.45
2:K:59:TYR:CE1	2:K:69:ILE:HG12	2.53	0.44
1:A:79:LEU:HB2	4:A:602:HOH:O	2.17	0.44
2:K:15:GLY:HA2	2:K:84:SER:HA	2.00	0.44
1:A:77:SER:HB3	1:A:83:ILE:HG12	1.99	0.44
1:A:247:LEU:HD23	1:A:247:LEU:HA	1.91	0.44
2:C:71[B]:ARG:HE	2:C:71[B]:ARG:HB3	1.42	0.44
1:A:152:SER:O	1:A:187:PRO:HD2	2.18	0.43
1:B:77:SER:HB3	1:B:83:ILE:HG12	2.01	0.43
2:C:53:SER:HA	2:C:71[B]:ARG:NH1	2.34	0.43
2:K:46:GLU:CD	4:K:201:HOH:O	2.61	0.42
1:A:182:ARG:NH1	4:A:604:HOH:O	2.52	0.42
2:K:40:ALA:HB3	2:K:43:LYS:HB2	2.02	0.42
1:B:64:LEU:HB3	1:B:65:PRO:HD3	2.00	0.42
1:A:64:LEU:HB3	1:A:65:PRO:HD3	2.02	0.42
1:B:154:LEU:HD22	1:B:183:LEU:HD12	2.02	0.42
2:C:35:THR:HG23	2:C:50:ALA:HB2	2.01	0.42
2:C:36:TRP:O	2:C:48:VAL:HB	2.20	0.42
1:A:262:GLY:HA3	1:A:330:PHE:CZ	2.55	0.41
2:K:47:TRP:HE1	2:K:50:ALA:HB2	1.85	0.41
1:B:151:LEU:HD21	1:B:183:LEU:HD13	2.02	0.41
1:B:150:LEU:HD23	1:B:150:LEU:HA	1.90	0.41
2:C:69:ILE:CG2	2:C:80:LEU:HD13	2.51	0.41
1:A:80:ILE:N	4:A:602:HOH:O	2.44	0.41
1:A:278:GLU:OE1	1:A:282:ARG:NH1	2.47	0.41
2:K:11:CYS:O	2:C:10:GLY:HA3	2.21	0.41
1:B:152:SER:O	1:B:187:PRO:HD2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:6:GLU:OE2	2:C:116:GLY:HA3	2.21	0.40
1:A:370:GLN:O	1:A:374:LEU:HG	2.21	0.40
1:B:262:GLY:HA3	1:B:330:PHE:CE2	2.56	0.40
1:A:421:GLY:O	2:C:103:LEU:HD23	2.21	0.40
1:A:77:SER:HB3	1:A:83:ILE:CG1	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	444/445 (100%)	434 (98%)	10 (2%)	0	100	100
1	B	442/445 (99%)	432 (98%)	10 (2%)	0	100	100
2	C	124/127 (98%)	117 (94%)	7 (6%)	0	100	100
2	K	124/127 (98%)	124 (100%)	0	0	100	100
All	All	1134/1144 (99%)	1107 (98%)	27 (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	
1	A	368/373 (99%)	361 (98%)	7 (2%)	50	69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	368/373 (99%)	360 (98%)	8 (2%)	45	65
2	C	99/100 (99%)	97 (98%)	2 (2%)	48	67
2	K	99/100 (99%)	94 (95%)	5 (5%)	21	32
All	All	934/946 (99%)	912 (98%)	22 (2%)	43	62

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	LEU
1	A	119	GLU
1	A	248	LYS
1	A	255	ASP
1	A	258	LEU
1	A	381	LYS
1	A	409	LEU
1	B	45	ASN
1	B	59	SER
1	B	86	GLN
1	B	104	LEU
1	B	248	LYS
1	B	258	LEU
1	B	409	LEU
1	B	442	ARG
2	K	25	SER
2	K	27	SER
2	K	52	THR
2	K	88	GLU
2	K	95	CYS
2	C	102	THR
2	C	120	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	139	GLN
1	A	285	ASN
1	A	413	HIS
1	B	34	GLN
1	B	139	GLN

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Mol	Chain	Res	Type
1	B	198	GLN
1	B	285	ASN
1	B	413	HIS
2	K	3	GLN
2	K	73	ASN
2	K	81	GLN
2	C	3	GLN

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 [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	443/445 (99%)	0.71	33 (7%)	20 22	42, 73, 120, 172	3 (0%)
1	B	441/445 (99%)	0.71	34 (7%)	19 21	27, 73, 117, 167	3 (0%)
2	C	124/127 (97%)	0.89	13 (10%)	11 13	30, 72, 99, 162	2 (1%)
2	K	124/127 (97%)	1.46	34 (27%)	1 1	44, 74, 112, 149	2 (1%)
All	All	1132/1144 (98%)	0.81	114 (10%)	12 14	27, 73, 118, 172	10 (0%)

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	167[A]	HIS	5.9
1	B	317	ILE	5.8
1	B	319	PHE	5.5
2	K	50	ALA	5.1
1	B	12	PRO	5.0
1	A	409	LEU	4.6
2	K	17	SER	4.6
1	B	318	SER	4.5
2	C	94	TYR	4.5
1	A	317	ILE	4.5
1	B	26	PHE	4.5
1	B	20	LEU	4.2
1	B	409	LEU	4.1
1	B	165	TRP	3.9
1	A	79	LEU	3.9
2	K	18	LEU	3.9
1	A	78	PRO	3.9
2	K	94	TYR	3.8
1	B	452	SER	3.8
2	K	69	ILE	3.7
2	K	103	LEU	3.7

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Mol	Chain	Res	Type	RSRZ
2	C	95	CYS	3.6
2	C	71[A]	ARG	3.6
2	K	71[A]	ARG	3.5
2	K	95	CYS	3.5
1	A	165	TRP	3.4
2	K	122	MET	3.3
1	A	323	VAL	3.3
2	K	14	ALA	3.2
1	B	167[A]	HIS	3.2
1	A	319	PHE	3.1
1	B	60	LEU	3.1
1	A	453	TRP	3.0
2	K	124	SER	3.0
2	K	82	MET	3.0
1	B	79	LEU	3.0
1	B	283	GLY	3.0
1	A	142	LEU	2.9
2	K	53	SER	2.9
2	C	124	SER	2.9
2	K	85	LEU	2.9
1	B	323	VAL	2.9
2	K	123	VAL	2.8
2	K	76	ASN	2.8
1	A	321	MET	2.8
2	C	122	MET	2.8
2	K	9	GLY	2.7
2	K	84	SER	2.7
1	A	185	HIS	2.7
2	C	69	ILE	2.7
1	B	29	PHE	2.7
2	C	14	ALA	2.7
1	A	320	GLY	2.6
1	B	151	LEU	2.6
2	K	83[A]	ASN	2.6
1	B	153	TYR	2.6
1	B	374	LEU	2.5
1	B	109	ARG	2.5
1	A	80	ILE	2.5
1	B	254	ALA	2.5
1	B	78	PRO	2.5
2	K	49	ALA	2.5
2	K	70	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	100	LEU	2.4
2	K	13	LYS	2.4
1	B	424	LEU	2.4
2	K	116	GLY	2.4
2	K	78	VAL	2.4
2	C	8	GLY	2.4
1	A	38	THR	2.4
1	A	256	LYS	2.4
1	A	318	SER	2.4
2	K	67	PHE	2.4
2	C	85	LEU	2.4
1	A	169	PHE	2.3
1	A	322	GLY	2.3
1	B	95	VAL	2.3
1	B	93	LEU	2.3
2	C	20	LEU	2.3
1	A	53	PRO	2.3
1	A	382	LEU	2.2
1	B	23	VAL	2.3
2	K	54	GLY	2.2
1	A	194	THR	2.2
1	A	54	THR	2.2
2	C	24	ALA	2.2
1	B	451	SER	2.2
2	K	93	TYR	2.2
2	K	11	CYS	2.2
1	A	82	LEU	2.2
1	B	322	GLY	2.2
2	C	97	ALA	2.2
2	K	10	GLY	2.1
1	A	106	ALA	2.1
1	B	125	ILE	2.1
1	B	174	LEU	2.1
2	K	22	CYS	2.1
1	B	123	THR	2.1
2	K	91	ALA	2.1
1	A	43	LYS	2.1
2	K	24	ALA	2.1
1	A	424	LEU	2.1
2	K	51	ILE	2.1
1	B	146	VAL	2.1
2	K	12	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	54	THR	2.0
1	A	104	LEU	2.0
1	A	145	LEU	2.0
2	C	76	ASN	2.0
1	B	72	ILE	2.0
1	A	26	PHE	2.0
1	B	330	PHE	2.0
1	A	149	HIS	2.0
1	A	12	PRO	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
3	ZN	A	501	1/1	1.00	0.01	69,69,69,69	0
3	ZN	B	501	1/1	1.00	0.03	61,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.