



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:26 PM UTC

PDB ID : 5F9F / pdb_00005f9f
Title : Crystal structure of RIG-I helicase-RD in complex with 24-mer blunt-end hairpin RNA
Authors : Wang, C.; Marcotrigiano, J.; Miller, M.T.; Jiang, F.
Deposited on : 2015-12-09
Resolution : 2.60 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

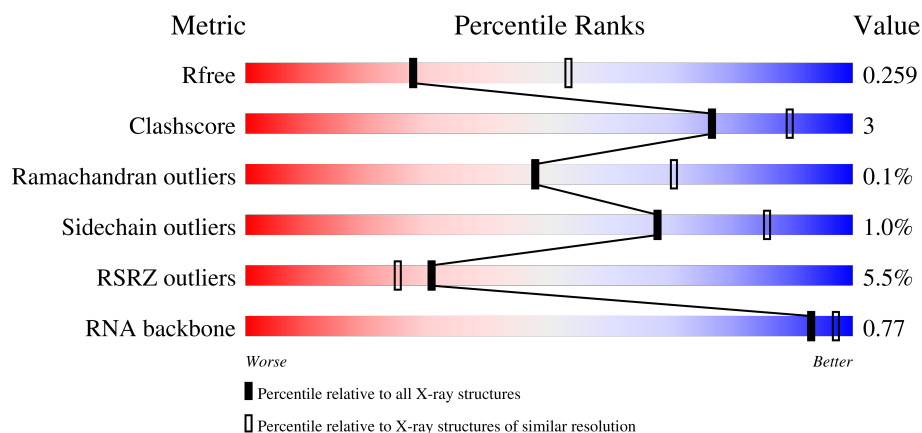
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)
RNA backbone	3983	1014 (2.84-2.36)

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	695	4% 90% 7%
1	C	695	5% 89% 7%
1	E	695	9% 88% 8%
1	G	695	4% 87% 8%

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Mol	Chain	Length	Quality of chain
1	I	695	
1	K	695	
2	B	24	
2	D	24	
2	F	24	
2	H	24	
2	J	24	
2	L	24	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 34918 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	673	Total	C	N	O	S	0	0	0
			5310	3393	902	982	33			
1	C	668	Total	C	N	O	S	0	0	0
			5266	3368	894	971	33			
1	E	672	Total	C	N	O	S	0	0	0
			5222	3338	882	969	33			
1	G	669	Total	C	N	O	S	0	0	0
			5282	3375	902	972	33			
1	I	674	Total	C	N	O	S	0	0	0
			5306	3395	899	979	33			
1	K	672	Total	C	N	O	S	0	0	0
			5256	3362	889	971	34			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	231	SER	-	expression tag	UNP O95786
C	231	SER	-	expression tag	UNP O95786
E	231	SER	-	expression tag	UNP O95786
G	231	SER	-	expression tag	UNP O95786
I	231	SER	-	expression tag	UNP O95786
K	231	SER	-	expression tag	UNP O95786

- Molecule 2 is a RNA chain called RNA (5'-R(*GP*AP*AP*UP*AP*UP*AP*AP*UP*AP*GP*UP*GP*AP*UP*AP*UP*UP*AP*UP*AP*UP*UP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	24	Total	C	N	O	P	0	0	0
			506	229	88	166	23			
2	D	24	Total	C	N	O	P	0	0	0
			506	229	88	166	23			
2	F	24	Total	C	N	O	P	0	0	0
			506	229	88	166	23			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	24	Total 506	C 229	N 88	O 166	P 23	0	0	0
2	J	24	Total 506	C 229	N 88	O 166	P 23	0	0	0
2	L	24	Total 506	C 229	N 88	O 166	P 23	0	0	0

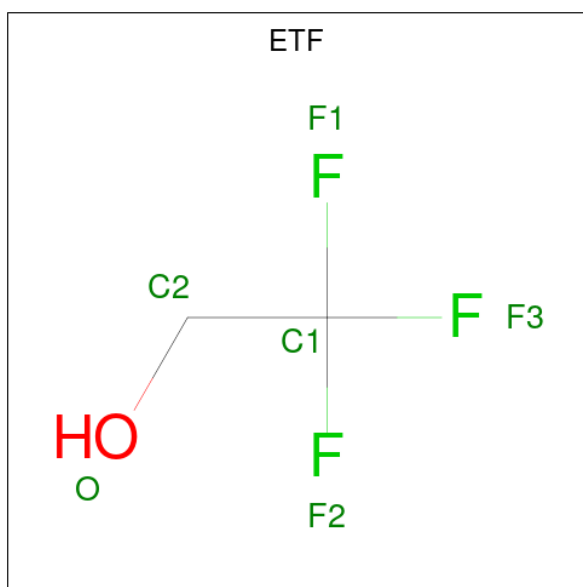
- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Zn 1	0	0
3	C	1	Total 1	Zn 1	0	0
3	E	1	Total 1	Zn 1	0	0
3	G	1	Total 1	Zn 1	0	0
3	I	1	Total 1	Zn 1	0	0
3	K	1	Total 1	Zn 1	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

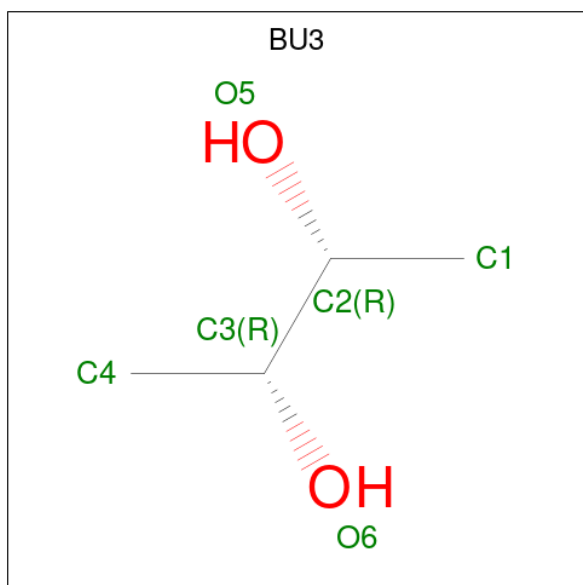
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Mg 1	0	0
4	C	1	Total 1	Mg 1	0	0
4	H	1	Total 1	Mg 1	0	0
4	I	2	Total 2	Mg 2	0	0
4	K	1	Total 1	Mg 1	0	0

- Molecule 5 is TRIFLUOROETHANOL (CCD ID: ETF) (formula: C₂H₃F₃O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	F	O	0	0
			6	2	3	1		
5	E	1	Total	C	F	O	0	0
			6	2	3	1		
5	G	1	Total	C	F	O	0	0
			6	2	3	1		
5	K	1	Total	C	F	O	0	0
			6	2	3	1		

- Molecule 6 is (R,R)-2,3-BUTANEDIOL (CCD ID: BU3) (formula: $C_4H_{10}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 6 4 2	0	0
6	A	1	Total C O 6 4 2	0	0
6	C	1	Total C O 6 4 2	0	0
6	C	1	Total C O 6 4 2	0	0
6	C	1	Total C O 6 4 2	0	0
6	C	1	Total C O 6 4 2	0	0
6	E	1	Total C O 6 4 2	0	0
6	E	1	Total C O 6 4 2	0	0
6	G	1	Total C O 6 4 2	0	0
6	G	1	Total C O 6 4 2	0	0
6	G	1	Total C O 6 4 2	0	0
6	I	1	Total C O 6 4 2	0	0
6	K	1	Total C O 6 4 2	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	34	Total O 34 34	0	0
7	B	2	Total O 2 2	0	0
7	C	14	Total O 14 14	0	0
7	D	4	Total O 4 4	0	0
7	E	16	Total O 16 16	0	0
7	F	3	Total O 3 3	0	0
7	G	21	Total O 21 21	0	0

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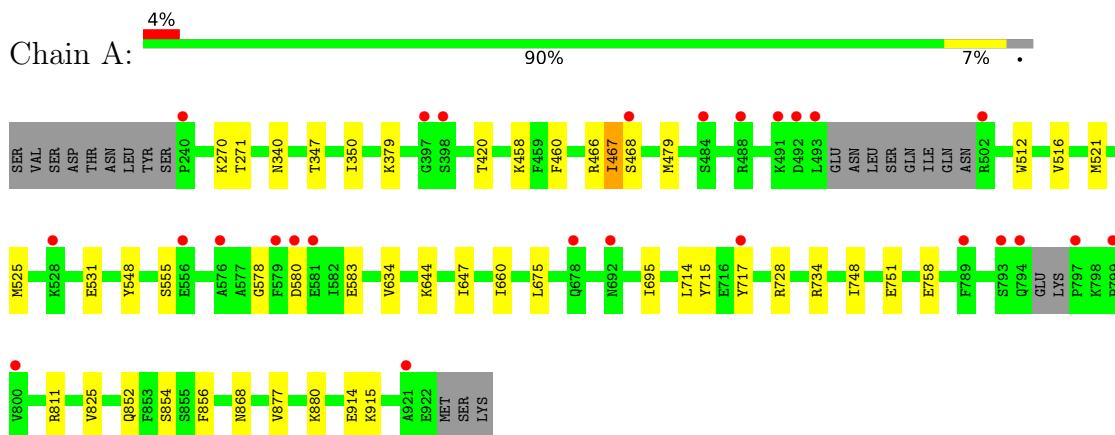
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	H	3	Total 3	O 3	0	0
7	I	7	Total 7	O 7	0	0
7	J	1	Total 1	O 1	0	0
7	K	20	Total 20	O 20	0	0
7	L	1	Total 1	O 1	0	0

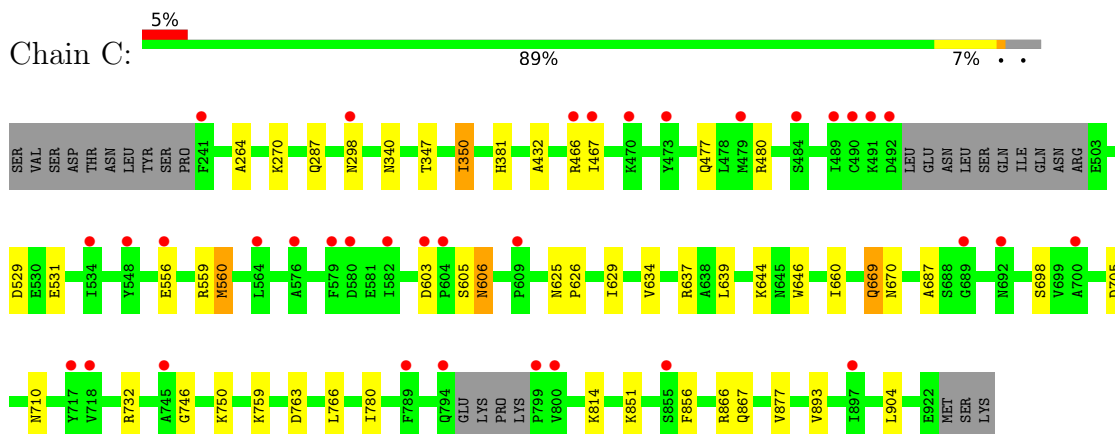
3 Residue-property plots [i](#)

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.

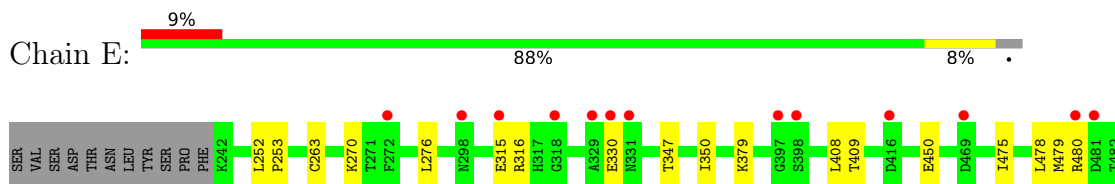
- Molecule 1: Probable ATP-dependent RNA helicase DDX58

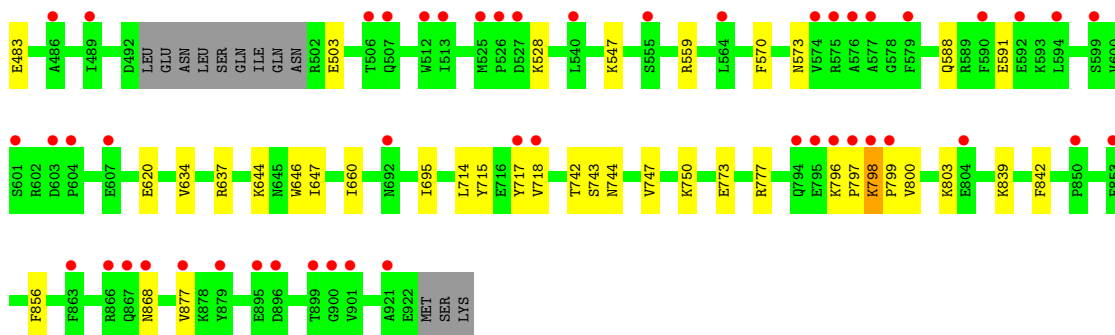


- Molecule 1: Probable ATP-dependent RNA helicase DDX58

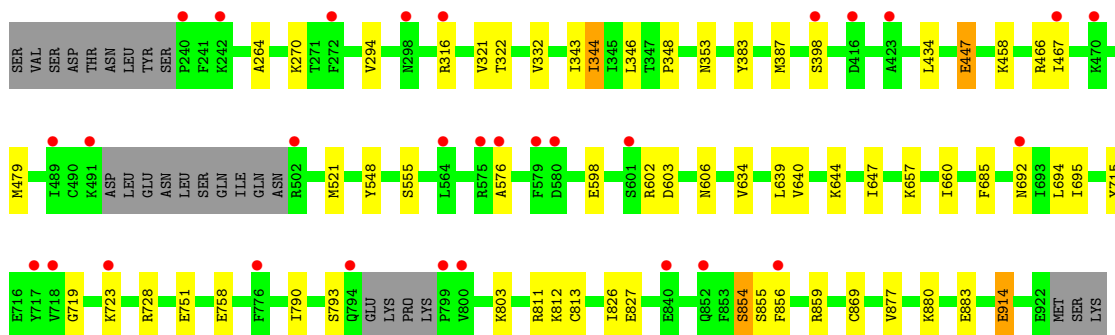
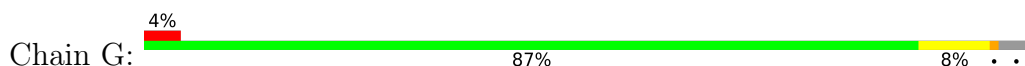


- Molecule 1: Probable ATP-dependent RNA helicase DDX58

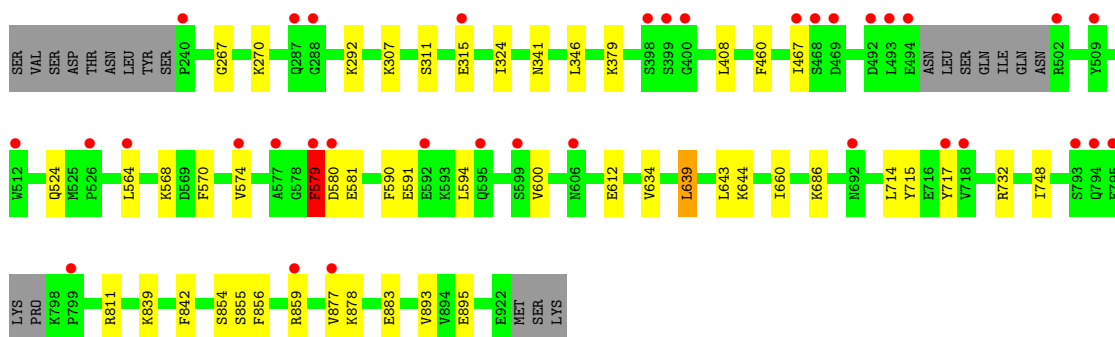
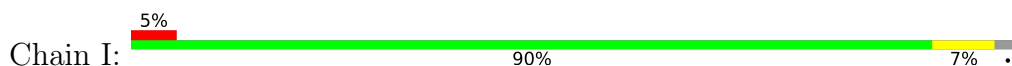




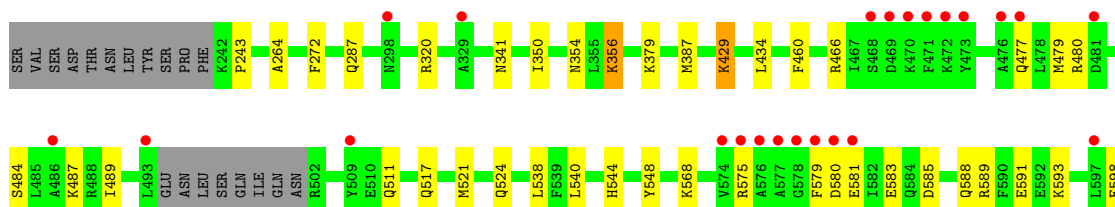
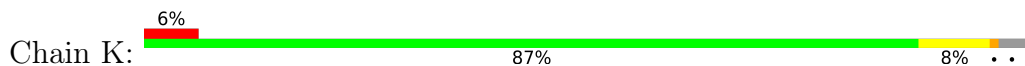
- Molecule 1: Probable ATP-dependent RNA helicase DDX58



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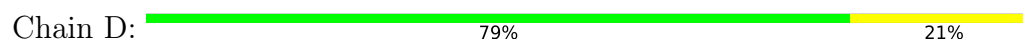




- Molecule 2: RNA (5'-R(*GP*AP*AP*UP*AP*UP*AP*AP*UP*AP*GP*UP*GP*AP*UP*AP*UP*UP*AP*UP*AP*UP*UP*C)-3')



- Molecule 2: RNA (5'-R(*GP*AP*AP*UP*AP*UP*AP*AP*UP*AP*GP*UP*GP*AP*UP*AP*UP*UP*AP*UP*AP*UP*UP*C)-3')



- Molecule 2: RNA (5'-R(*GP*AP*AP*UP*AP*UP*AP*AP*UP*AP*GP*UP*GP*AP*UP*AP*UP*UP*AP*UP*AP*UP*UP*C)-3')



- Molecule 2: RNA (5'-R(*GP*AP*AP*UP*AP*UP*AP*AP*UP*AP*GP*UP*GP*AP*UP*AP*UP*UP*AP*UP*AP*UP*UP*C)-3')

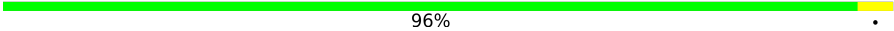


There are no outlier residues recorded for this chain.

- Molecule 2: RNA (5'-R(*GP*AP*AP*UP*AP*UP*AP*AP*UP*AP*GP*UP*GP*AP*UP*AP*UP*UP*AP*UP*AP*UP*UP*C)-3')



- Molecule 2: RNA (5'-R(*GP*AP*AP*UP*AP*UP*AP*AP*UP*AP*GP*UP*GP*AP*UP*AP*UP*UP*AP*UP*AP*UP*UP*C)-3')

Chain L:  96%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	111.52Å 174.25Å 308.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.99 – 2.60 29.99 – 2.60	Depositor EDS
% Data completeness (in resolution range)	90.2 (29.99-2.60) 87.2 (29.99-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.9_1690	Depositor
R, R_{free}	0.203 , 0.259 0.206 , 0.259	Depositor DCC
R_{free} test set	8318 reflections (4.51%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.323	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 50.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	34918	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.73% 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: ETF, MG, ZN, BU3

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	0.27	0/5419	0.71	2/7325 (0.0%)
1	C	0.27	0/5373	0.70	0/7260
1	E	0.29	0/5330	0.73	1/7222 (0.0%)
1	G	0.28	0/5390	0.70	1/7284 (0.0%)
1	I	0.29	0/5414	0.72	3/7315 (0.0%)
1	K	0.28	0/5362	0.71	4/7251 (0.1%)
2	B	0.06	0/566	0.15	0/879
2	D	0.06	0/566	0.15	0/879
2	F	0.06	0/566	0.14	0/879
2	H	0.06	0/566	0.15	0/879
2	J	0.06	0/566	0.15	0/879
2	L	0.06	0/566	0.15	0/879
All	All	0.27	0/35684	0.68	11/48931 (0.0%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	854	SER	CB-CA-C	-5.62	110.08	116.54
1	K	607	GLU	N-CA-C	5.56	116.97	108.52
1	A	854	SER	CB-CA-C	-5.52	110.19	116.54
1	G	854	SER	CB-CA-C	-5.45	110.27	116.54
1	I	379	LYS	CB-CA-C	-5.29	110.46	116.54
1	A	379	LYS	CB-CA-C	-5.26	110.48	116.54
1	E	379	LYS	CB-CA-C	-5.22	110.54	116.54
1	K	884	ILE	N-CA-C	5.20	113.61	107.84
1	K	379	LYS	CB-CA-C	-5.16	110.61	116.54
1	I	579	PHE	N-CA-C	5.11	117.57	109.24
1	I	600	VAL	N-CA-C	-5.01	105.51	110.62

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	5310	0	5266	26	0
1	C	5266	0	5231	33	0
1	E	5222	0	5107	33	0
1	G	5282	0	5242	36	0
1	I	5306	0	5267	31	0
1	K	5256	0	5183	39	0
2	B	506	0	256	1	0
2	D	506	0	256	3	0
2	F	506	0	256	1	0
2	H	506	0	256	0	0
2	J	506	0	256	1	0
2	L	506	0	256	1	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
3	G	1	0	0	0	0
3	I	1	0	0	0	0
3	K	1	0	0	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
4	H	1	0	0	0	0
4	I	2	0	0	0	0
4	K	1	0	0	0	0
5	A	6	0	3	1	0
5	E	6	0	3	0	0
5	G	6	0	3	1	0
5	K	6	0	3	0	0
6	A	12	0	20	1	0
6	C	24	0	40	4	0
6	E	12	0	20	4	0
6	G	18	0	30	5	0
6	I	6	0	10	3	0
6	K	6	0	10	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	34	0	0	0	0
7	B	2	0	0	0	0
7	C	14	0	0	0	0
7	D	4	0	0	0	0
7	E	16	0	0	0	0
7	F	3	0	0	0	0
7	G	21	0	0	0	0
7	H	3	0	0	0	0
7	I	7	0	0	0	0
7	J	1	0	0	0	0
7	K	20	0	0	0	0
7	L	1	0	0	0	0
All	All	34918	0	32974	193	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 3.

All (193) 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:K:600:VAL:O	1:K:606:ASN:ND2	2.03	0.92
1:G:826:ILE:HG22	1:G:827:GLU:HG3	1.67	0.76
1:E:480:ARG:HA	1:E:483:GLU:HG2	1.73	0.70
1:G:880:LYS:O	1:K:287:GLN:NE2	2.19	0.69
1:I:570:PHE:O	1:I:574:VAL:HG23	1.93	0.68
1:K:603:ASP:HB3	1:K:606:ASN:ND2	2.09	0.68
1:I:580:ASP:OD1	1:I:581:GLU:N	2.27	0.67
1:A:467:ILE:HD12	1:A:468:SER:H	1.58	0.67
1:K:580:ASP:OD1	1:K:581:GLU:N	2.28	0.66
1:A:466:ARG:NH1	1:A:555:SER:O	2.30	0.65
1:G:458:LYS:NZ	1:G:751:GLU:OE2	2.27	0.64
1:C:298:ASN:HD21	1:C:381:HIS:CE1	2.15	0.64
1:C:477:GLN:HA	1:C:480:ARG:HG3	1.79	0.64
1:K:603:ASP:HB3	1:K:606:ASN:HD21	1.64	0.62
1:I:270:LYS:HB3	6:I:1004:BU3:H2	1.81	0.62
1:G:723:LYS:HG2	6:G:1005:BU3:H42	1.82	0.61
1:G:657:LYS:HD2	1:G:692:ASN:HB2	1.83	0.61
1:G:803:LYS:HG2	1:G:914:GLU:HG3	1.83	0.61
1:G:644:LYS:HD2	1:G:660:ILE:HD11	1.83	0.60
1:K:477:GLN:HA	1:K:480:ARG:HB2	1.84	0.60
1:I:856:PHE:HD2	1:I:877:VAL:HG21	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:603:ASP:HB3	1:C:606:ASN:OD1	2.01	0.59
1:E:798:LYS:H	1:E:799:PRO:CD	2.15	0.58
1:C:264:ALA:HB3	6:C:1006:BU3:H2	1.83	0.58
1:E:717:TYR:HE2	6:E:1004:BU3:H2	1.69	0.58
1:A:644:LYS:HD2	1:A:660:ILE:HD11	1.87	0.57
1:K:489:ILE:HD11	1:K:538:LEU:HG	1.87	0.57
1:C:759:LYS:NZ	1:C:763:ASP:OD2	2.37	0.56
1:C:669:GLN:OE1	1:C:670:ASN:N	2.35	0.56
1:C:866:ARG:HG2	1:C:867:GLN:HG3	1.86	0.56
1:E:798:LYS:H	1:E:799:PRO:HD3	1.69	0.56
1:A:880:LYS:O	1:C:287:GLN:NE2	2.35	0.56
1:C:746:GLY:O	1:C:750:LYS:HG3	2.06	0.56
1:A:271:THR:HB	6:A:1005:BU3:H13	1.88	0.55
1:A:479:MET:HE2	1:A:548:TYR:HB3	1.89	0.55
1:E:475:ILE:O	1:E:479:MET:HG3	2.07	0.55
1:G:603:ASP:OD2	1:G:606:ASN:ND2	2.40	0.54
1:K:511:GLN:NE2	2:L:9:U:O2	2.38	0.54
1:C:340:ASN:HB2	6:C:1004:BU3:H43	1.90	0.53
1:C:644:LYS:HD2	1:C:660:ILE:HD11	1.90	0.53
1:K:575:ARG:HA	1:K:579:PHE:HE1	1.74	0.53
1:I:714:LEU:HB3	1:I:717:TYR:HB3	1.91	0.53
1:A:734:ARG:NH1	1:E:450:GLU:OE2	2.31	0.53
1:I:859:ARG:HH12	1:I:878:LYS:N	2.08	0.52
1:G:719:GLY:HA3	6:G:1005:BU3:H41	1.91	0.52
1:K:724:MET:HE1	1:K:755:MET:HG3	1.92	0.52
1:E:570:PHE:O	1:E:573:ASN:HB2	2.09	0.52
1:C:603:ASP:OD1	1:C:605:SER:N	2.35	0.51
1:G:294:VAL:HG23	1:G:346:LEU:HD23	1.91	0.51
1:E:714:LEU:HB3	1:E:717:TYR:HB3	1.92	0.51
1:K:517:GLN:HG2	1:K:521:MET:HE3	1.93	0.51
1:K:607:GLU:HA	1:K:607:GLU:OE1	2.11	0.51
1:I:590:PHE:CE2	1:I:594:LEU:HD13	2.47	0.50
1:I:564:LEU:HG	1:I:568:LYS:HE3	1.93	0.50
1:E:773:GLU:OE2	1:E:777:ARG:NH2	2.45	0.50
1:I:859:ARG:HH12	1:I:877:VAL:C	2.19	0.50
1:A:512:TRP:CZ2	1:A:516:VAL:HG21	2.47	0.50
1:I:270:LYS:CB	6:I:1004:BU3:H2	2.41	0.49
1:I:639:LEU:HD23	1:I:643:LEU:HG	1.94	0.49
1:E:647:ILE:HD12	1:E:695:ILE:HD11	1.94	0.49
1:C:893:VAL:HG22	1:C:904:LEU:HD11	1.95	0.48
1:I:574:VAL:HG12	1:I:579:PHE:HE1	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:637:ARG:NH1	2:D:21:A:OP2	2.45	0.48
1:G:447:GLU:H	1:G:447:GLU:CD	2.21	0.48
1:I:859:ARG:HH21	1:I:883:GLU:HB3	1.78	0.48
1:G:598:GLU:HG2	1:G:602:ARG:CZ	2.43	0.48
1:I:839:LYS:HA	1:I:842:PHE:CE2	2.48	0.48
2:D:11:G:N2	2:D:14:A:OP2	2.41	0.48
1:E:796:LYS:N	1:E:797:PRO:HD3	2.29	0.48
1:G:264:ALA:HB3	6:G:1003:BU3:H43	1.95	0.48
1:I:644:LYS:HD2	1:I:660:ILE:HD11	1.95	0.48
1:K:759:LYS:NZ	1:K:763:ASP:OD2	2.45	0.48
1:K:350:ILE:O	1:K:354:ASN:ND2	2.46	0.47
1:A:825:VAL:HG21	1:A:915:LYS:HB3	1.94	0.47
1:K:803:LYS:HE3	1:K:803:LYS:HB3	1.62	0.47
1:K:860:ALA:O	1:K:876:HIS:N	2.35	0.47
1:E:408:LEU:HD22	6:E:1003:BU3:H41	1.95	0.47
2:J:11:G:N2	2:J:14:A:OP2	2.40	0.47
1:C:705:ASP:CG	1:C:732:ARG:HH22	2.22	0.47
1:G:322:THR:HG23	1:G:344:ILE:HG23	1.95	0.47
1:K:264:ALA:HB3	6:K:1004:BU3:H41	1.96	0.47
1:G:859:ARG:NE	1:G:883:GLU:OE2	2.42	0.47
1:C:477:GLN:HA	1:C:480:ARG:CG	2.44	0.47
1:A:420:THR:OG1	1:E:620:GLU:OE1	2.33	0.47
1:A:458:LYS:NZ	1:A:751:GLU:OE2	2.44	0.47
1:C:867:GLN:HG2	1:G:576:ALA:HA	1.95	0.47
1:G:856:PHE:HD2	1:G:877:VAL:HG21	1.78	0.47
1:I:311:SER:O	1:I:315:GLU:HG3	2.13	0.47
1:G:854:SER:OG	1:G:855:SER:N	2.42	0.47
1:K:540:LEU:O	1:K:544:HIS:ND1	2.48	0.47
1:E:263:CYS:HA	1:E:409:THR:O	2.15	0.47
1:G:790:ILE:O	1:G:793:SER:OG	2.33	0.47
1:C:634:VAL:HB	1:C:639:LEU:HD23	1.97	0.46
1:E:588:GLN:O	1:E:591:GLU:N	2.48	0.46
1:A:852:GLN:HG3	1:A:856:PHE:O	2.15	0.46
1:C:698:SER:HG	2:D:20:U:HO2'	1.62	0.46
1:E:718:VAL:HG13	1:E:750:LYS:HG2	1.98	0.46
1:I:856:PHE:CD2	1:I:877:VAL:HG21	2.50	0.46
1:C:477:GLN:O	1:C:477:GLN:HG3	2.14	0.46
1:A:578:GLY:HA3	1:A:868:ASN:ND2	2.30	0.46
1:E:270:LYS:HB3	6:E:1003:BU3:H2	1.98	0.46
1:G:321:VAL:HG22	1:G:343:ILE:HB	1.98	0.46
1:E:839:LYS:HA	1:E:842:PHE:CE2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:429:LYS:HB3	1:K:429:LYS:HE3	1.64	0.45
1:A:714:LEU:HB3	1:A:717:TYR:HB3	1.98	0.45
1:A:634:VAL:HG12	1:A:715:TYR:HB3	1.98	0.45
1:A:340:ASN:HB2	6:C:1004:BU3:H3	1.99	0.45
1:A:347:THR:HB	1:A:350:ILE:HD13	1.98	0.45
1:E:347:THR:HB	1:E:350:ILE:HD13	1.99	0.45
1:I:591:GLU:HA	1:I:594:LEU:HB2	1.99	0.45
1:K:588:GLN:HA	1:K:591:GLU:HB2	1.98	0.45
1:K:606:ASN:O	1:K:607:GLU:HB2	2.17	0.45
1:C:856:PHE:HD2	1:C:877:VAL:HG21	1.81	0.45
1:I:408:LEU:HB3	6:I:1004:BU3:H13	1.98	0.44
1:G:634:VAL:HG12	1:G:715:TYR:HB3	1.98	0.44
1:E:742:THR:OG1	1:E:743:SER:N	2.51	0.44
1:K:568:LYS:NZ	1:K:598:GLU:OE2	2.25	0.44
1:A:460:PHE:HB3	1:A:748:ILE:HD12	1.99	0.44
1:G:728:ARG:NH2	1:G:758:GLU:OE1	2.50	0.44
1:E:528:LYS:HA	1:E:528:LYS:HD2	1.65	0.44
1:G:466:ARG:NH1	1:G:555:SER:O	2.51	0.44
1:K:634:VAL:HB	1:K:639:LEU:HD23	1.99	0.44
1:G:332:VAL:HG11	1:K:320:ARG:HD2	2.00	0.44
1:K:243:PRO:HB3	1:K:272:PHE:HE2	1.83	0.43
1:K:585:ASP:O	1:K:589:ARG:HG3	2.18	0.43
1:G:812:LYS:HD3	1:G:869:CYS:SG	2.59	0.43
1:C:531:GLU:OE1	1:C:814:LYS:NZ	2.45	0.43
1:E:252:LEU:HB3	1:E:253:PRO:HD3	1.99	0.43
1:K:580:ASP:HB3	1:K:583:GLU:HB2	2.00	0.43
1:A:647:ILE:HD12	1:A:695:ILE:HD11	2.00	0.43
1:C:270:LYS:HD3	6:C:1006:BU3:H43	2.00	0.43
1:G:387:MET:HE2	1:G:434:LEU:HA	2.00	0.43
1:I:859:ARG:HH12	1:I:877:VAL:HA	1.84	0.43
1:E:634:VAL:HG12	1:E:715:TYR:HB3	2.01	0.43
1:I:574:VAL:HG12	1:I:579:PHE:CE1	2.54	0.43
1:I:854:SER:OG	1:I:855:SER:N	2.52	0.43
1:E:559:ARG:HD3	1:E:646:TRP:HD1	1.84	0.43
1:A:270:LYS:HB3	5:A:1003:ETF:H21	2.01	0.42
1:E:717:TYR:CE2	6:E:1004:BU3:H2	2.50	0.42
1:E:637:ARG:NH1	2:F:21:A:OP2	2.49	0.42
1:E:644:LYS:HD2	1:E:660:ILE:HD11	1.99	0.42
1:K:484:SER:HA	1:K:487:LYS:HG3	2.00	0.42
1:C:560:MET:HB2	1:C:560:MET:HE3	1.58	0.42
1:G:634:VAL:HB	1:G:639:LEU:HD23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:859:ARG:HH12	1:I:877:VAL:CA	2.32	0.42
1:A:347:THR:HG21	2:B:23:U:O3'	2.20	0.42
1:A:580:ASP:HB3	1:A:583:GLU:HB2	2.01	0.42
1:C:529:ASP:OD1	1:C:529:ASP:N	2.53	0.42
1:E:547:LYS:HG3	1:E:570:PHE:CE1	2.55	0.42
1:I:324:ILE:HB	1:I:346:LEU:HD12	2.01	0.42
1:K:387:MET:HE2	1:K:434:LEU:HA	2.02	0.42
1:A:525:MET:HB2	1:A:531:GLU:HB2	2.02	0.41
1:C:629:ILE:HD11	1:C:687:ALA:HA	2.00	0.41
1:I:634:VAL:HG12	1:I:715:TYR:HB3	2.00	0.41
1:K:856:PHE:HD2	1:K:877:VAL:HG21	1.85	0.41
1:E:315:GLU:OE2	1:I:307:LYS:NZ	2.53	0.41
1:E:856:PHE:HD2	1:E:877:VAL:HG21	1.86	0.41
1:E:868:ASN:C	1:E:868:ASN:OD1	2.62	0.41
1:K:356:LYS:HE2	1:K:356:LYS:HB3	1.79	0.41
1:A:728:ARG:NH2	1:A:758:GLU:OE1	2.52	0.41
1:C:466:ARG:HH11	1:C:466:ARG:HG2	1.84	0.41
1:G:353:ASN:HD22	6:G:1004:BU3:H11	1.85	0.41
1:G:640:VAL:HG12	1:G:660:ILE:HD12	2.01	0.41
1:K:466:ARG:NH2	1:K:607:GLU:HA	2.35	0.41
1:E:744:ASN:HB3	1:E:747:VAL:HG23	2.01	0.41
1:G:521:MET:SD	1:G:811:ARG:HD3	2.61	0.41
1:G:647:ILE:HD12	1:G:695:ILE:HD11	2.01	0.41
1:G:803:LYS:HG2	1:G:914:GLU:CG	2.48	0.41
1:C:347:THR:HB	1:C:350:ILE:HD13	2.02	0.41
1:I:811:ARG:HG2	1:I:893:VAL:HG23	2.02	0.41
1:K:664:ARG:O	1:K:849:LYS:NZ	2.43	0.41
1:G:479:MET:HE2	1:G:548:TYR:HB3	2.02	0.41
1:K:479:MET:HE2	1:K:548:TYR:HB3	2.02	0.41
1:A:521:MET:SD	1:A:811:ARG:HD3	2.61	0.41
1:C:559:ARG:HD3	1:C:646:TRP:HD1	1.86	0.41
1:G:270:LYS:HB3	6:G:1003:BU3:H2	2.03	0.41
1:G:685:PHE:CE1	1:G:694:LEU:HB2	2.56	0.41
1:I:460:PHE:HB3	1:I:748:ILE:HD12	2.03	0.41
1:K:460:PHE:HB3	1:K:748:ILE:HD12	2.03	0.41
1:K:854:SER:OG	1:K:855:SER:N	2.50	0.41
1:C:629:ILE:H	1:C:710:ASN:HD21	1.69	0.40
1:K:466:ARG:HH21	1:K:607:GLU:HA	1.86	0.40
1:C:432:ALA:HA	1:C:780:ILE:HG23	2.03	0.40
1:E:478:LEU:HA	1:E:478:LEU:HD23	1.83	0.40
1:A:856:PHE:HD2	1:A:877:VAL:HG21	1.87	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:298:ASN:HD21	1:C:381:HIS:HE1	1.64	0.40
1:C:625:ASN:HA	1:C:626:PRO:HD2	1.98	0.40
1:I:267:GLY:HA3	1:I:732:ARG:HD3	2.04	0.40
1:I:292:LYS:NZ	1:I:341:ASN:O	2.48	0.40
1:K:664:ARG:HH11	1:K:675:LEU:HD12	1.86	0.40
1:G:348:PRO:HG3	1:G:383:TYR:CE1	2.56	0.40
5:G:1002:ETF:H21	1:K:341:ASN:OD1	2.22	0.40
1:I:524:GLN:NE2	1:I:895:GLU:OE2	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	667/695 (96%)	642 (96%)	24 (4%)	1 (0%)	48	70
1	C	662/695 (95%)	633 (96%)	29 (4%)	0	100	100
1	E	668/695 (96%)	628 (94%)	39 (6%)	1 (0%)	48	70
1	G	663/695 (95%)	635 (96%)	28 (4%)	0	100	100
1	I	668/695 (96%)	635 (95%)	33 (5%)	0	100	100
1	K	666/695 (96%)	635 (95%)	31 (5%)	0	100	100
All	All	3994/4170 (96%)	3808 (95%)	184 (5%)	2 (0%)	48	70

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	798	LYS
1	A	914	GLU

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	578/623 (93%)	576 (100%)	2 (0%)	86	94
1	C	572/623 (92%)	564 (99%)	8 (1%)	59	81
1	E	558/623 (90%)	552 (99%)	6 (1%)	65	84
1	G	573/623 (92%)	566 (99%)	7 (1%)	63	83
1	I	575/623 (92%)	570 (99%)	5 (1%)	70	87
1	K	565/623 (91%)	558 (99%)	7 (1%)	63	83
All	All	3421/3738 (92%)	3386 (99%)	35 (1%)	68	86

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	467	ILE
1	A	675	LEU
1	C	350	ILE
1	C	467	ILE
1	C	556	GLU
1	C	560	MET
1	C	606	ASN
1	C	669	GLN
1	C	766	LEU
1	C	851	LYS
1	E	276	LEU
1	E	316	ARG
1	E	330	GLU
1	E	503	GLU
1	E	800	VAL
1	E	803	LYS
1	G	316	ARG
1	G	344	ILE
1	G	398	SER
1	G	447	GLU
1	G	467	ILE
1	G	813	CYS

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Mol	Chain	Res	Type
1	G	914	GLU
1	I	467	ILE
1	I	579	PHE
1	I	612	GLU
1	I	639	LEU
1	I	686	LYS
1	K	356	LYS
1	K	429	LYS
1	K	524	GLN
1	K	593	LYS
1	K	607	GLU
1	K	803	LYS
1	K	840	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	289	GLN
1	A	298	ASN
1	A	317	HIS
1	A	619	GLN
1	C	298	ASN
1	C	299	GLN
1	C	305	GLN
1	C	340	ASN
1	C	349	GLN
1	C	389	ASN
1	C	507	GLN
1	C	573	ASN
1	E	340	ASN
1	E	349	GLN
1	E	389	ASN
1	G	298	ASN
1	G	305	GLN
1	G	306	GLN
1	G	317	HIS
1	G	389	ASN
1	G	606	ASN
1	G	868	ASN
1	I	305	GLN
1	I	306	GLN
1	I	511	GLN

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Mol	Chain	Res	Type
1	I	549	ASN
1	I	619	GLN
1	I	782	HIS
1	K	298	ASN
1	K	306	GLN
1	K	349	GLN
1	K	393	GLN
1	K	517	GLN
1	K	524	GLN
1	K	606	ASN
1	K	691	HIS
1	K	752	GLN
1	K	871	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	23/24 (95%)	0	0
2	D	23/24 (95%)	1 (4%)	0
2	F	23/24 (95%)	0	0
2	H	23/24 (95%)	0	0
2	J	23/24 (95%)	0	0
2	L	23/24 (95%)	0	0
All	All	138/144 (95%)	1 (0%)	0

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	D	15	U

There are no RNA pucker outliers to report.

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 29 ligands modelled in this entry, 12 are monoatomic - leaving 17 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	ETF	K	1003	-	5,5,5	0.41	0	7,7,7	0.65	0
6	BU3	E	1004	-	4,5,5	0.56	0	6,6,6	0.42	0
6	BU3	G	1005	-	4,5,5	0.51	0	6,6,6	0.50	0
6	BU3	A	1005	-	4,5,5	0.50	0	6,6,6	0.37	0
6	BU3	K	1004	-	4,5,5	0.51	0	6,6,6	0.39	0
6	BU3	C	1003	-	4,5,5	0.50	0	6,6,6	0.39	0
6	BU3	G	1004	-	4,5,5	0.51	0	6,6,6	0.41	0
6	BU3	C	1006	-	4,5,5	0.49	0	6,6,6	0.38	0
6	BU3	G	1003	-	4,5,5	0.53	0	6,6,6	0.36	0
6	BU3	C	1005	-	4,5,5	0.52	0	6,6,6	0.41	0
5	ETF	E	1002	-	5,5,5	0.39	0	7,7,7	0.65	0
6	BU3	E	1003	-	4,5,5	0.54	0	6,6,6	0.38	0
5	ETF	A	1003	-	5,5,5	0.40	0	7,7,7	0.65	0
6	BU3	I	1004	-	4,5,5	0.53	0	6,6,6	0.34	0
5	ETF	G	1002	-	5,5,5	0.39	0	7,7,7	0.68	0
6	BU3	A	1004	-	4,5,5	0.51	0	6,6,6	0.41	0
6	BU3	C	1004	-	4,5,5	0.55	0	6,6,6	0.37	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	ETF	K	1003	-	-	0/3/3/3	-
6	BU3	E	1004	-	-	4/4/4/4	-
6	BU3	G	1005	-	-	4/4/4/4	-
6	BU3	A	1005	-	-	4/4/4/4	-
6	BU3	K	1004	-	-	4/4/4/4	-
6	BU3	C	1003	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	BU3	G	1004	-	-	4/4/4/4	-
6	BU3	C	1006	-	-	4/4/4/4	-
6	BU3	G	1003	-	-	0/4/4/4	-
6	BU3	C	1005	-	-	0/4/4/4	-
5	ETF	E	1002	-	-	1/3/3/3	-
6	BU3	E	1003	-	-	2/4/4/4	-
5	ETF	A	1003	-	-	0/3/3/3	-
6	BU3	I	1004	-	-	0/4/4/4	-
5	ETF	G	1002	-	-	3/3/3/3	-
6	BU3	A	1004	-	-	2/4/4/4	-
6	BU3	C	1004	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1005	BU3	C1-C2-C3-O6
6	A	1005	BU3	O5-C2-C3-C4
6	A	1005	BU3	C1-C2-C3-C4
6	C	1004	BU3	C1-C2-C3-O6
6	C	1004	BU3	O5-C2-C3-C4
6	C	1004	BU3	C1-C2-C3-C4
6	C	1006	BU3	C1-C2-C3-O6
6	C	1006	BU3	O5-C2-C3-C4
6	C	1006	BU3	C1-C2-C3-C4
6	E	1004	BU3	O5-C2-C3-O6
6	E	1004	BU3	C1-C2-C3-C4
6	G	1004	BU3	O5-C2-C3-O6
6	G	1004	BU3	C1-C2-C3-O6
6	G	1004	BU3	O5-C2-C3-C4
6	G	1005	BU3	O5-C2-C3-O6
6	G	1005	BU3	C1-C2-C3-O6
6	G	1005	BU3	O5-C2-C3-C4
6	K	1004	BU3	C1-C2-C3-O6
6	K	1004	BU3	O5-C2-C3-C4
6	K	1004	BU3	C1-C2-C3-C4
6	A	1004	BU3	C1-C2-C3-O6

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Mol	Chain	Res	Type	Atoms
6	A	1004	BU3	O5-C2-C3-C4
6	E	1003	BU3	C1-C2-C3-O6
6	E	1003	BU3	O5-C2-C3-C4
6	E	1004	BU3	C1-C2-C3-O6
6	E	1004	BU3	O5-C2-C3-C4
6	G	1004	BU3	C1-C2-C3-C4
6	G	1005	BU3	C1-C2-C3-C4
6	A	1005	BU3	O5-C2-C3-O6
6	C	1004	BU3	O5-C2-C3-O6
6	C	1006	BU3	O5-C2-C3-O6
6	K	1004	BU3	O5-C2-C3-O6
5	G	1002	ETF	F2-C1-C2-O
5	G	1002	ETF	F1-C1-C2-O
5	G	1002	ETF	F3-C1-C2-O
5	E	1002	ETF	F2-C1-C2-O

There are no ring outliers.

12 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	E	1004	BU3	2	0
6	G	1005	BU3	2	0
6	A	1005	BU3	1	0
6	K	1004	BU3	1	0
6	G	1004	BU3	1	0
6	C	1006	BU3	2	0
6	G	1003	BU3	2	0
6	E	1003	BU3	2	0
5	A	1003	ETF	1	0
6	I	1004	BU3	3	0
5	G	1002	ETF	1	0
6	C	1004	BU3	2	0

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	673/695 (96%)	0.05	26 (3%)	43	38	17, 43, 85, 138	0
1	C	668/695 (96%)	0.27	35 (5%)	33	27	21, 50, 95, 150	0
1	E	672/695 (96%)	0.50	62 (9%)	14	11	24, 55, 107, 152	0
1	G	669/695 (96%)	0.26	30 (4%)	38	32	25, 50, 88, 124	0
1	I	674/695 (96%)	0.41	35 (5%)	33	27	27, 55, 96, 136	0
1	K	672/695 (96%)	0.37	39 (5%)	29	23	24, 55, 99, 146	0
2	B	24/24 (100%)	-0.17	1 (4%)	40	35	30, 38, 99, 142	0
2	D	24/24 (100%)	-0.19	0	100	100	25, 39, 121, 124	0
2	F	24/24 (100%)	-0.22	0	100	100	34, 43, 100, 109	0
2	H	24/24 (100%)	-0.29	0	100	100	31, 42, 109, 115	0
2	J	24/24 (100%)	-0.25	0	100	100	34, 46, 107, 110	0
2	L	24/24 (100%)	-0.27	0	100	100	34, 46, 111, 118	0
All	All	4172/4314 (96%)	0.29	228 (5%)	30	25	17, 51, 98, 152	0

All (228) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	717	TYR	5.6
1	A	797	PRO	5.6
1	A	579	PHE	5.0
1	C	470	LYS	4.9
1	C	492	ASP	4.8
1	A	493	LEU	4.8
1	K	796	LYS	4.6
1	A	576	ALA	4.3
1	C	700	ALA	4.3
1	E	318	GLY	4.2
1	I	494	GLU	4.2

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Mol	Chain	Res	Type	RSRZ
1	I	718	VAL	4.1
1	I	469	ASP	4.1
1	G	240	PRO	4.0
1	K	578	GLY	4.0
1	G	718	VAL	4.0
1	E	329	ALA	4.0
1	K	579	PHE	4.0
1	C	717	TYR	3.9
1	I	493	LEU	3.9
1	E	798	LYS	3.8
1	E	489	ILE	3.8
1	K	470	LYS	3.7
1	C	799	PRO	3.7
1	I	877	VAL	3.7
1	K	794	GLN	3.7
1	A	717	TYR	3.7
1	E	794	GLN	3.6
1	G	580	ASP	3.6
1	I	717	TYR	3.6
1	G	692	ASN	3.5
1	K	692	ASN	3.5
1	E	398	SER	3.5
1	G	470	LYS	3.5
1	C	579	PHE	3.5
1	K	608	ASN	3.4
1	I	579	PHE	3.4
1	A	240	PRO	3.4
1	C	604	PRO	3.4
1	E	900	GLY	3.4
1	C	603	ASP	3.4
1	E	718	VAL	3.4
1	E	526	PRO	3.3
1	G	502	ARG	3.3
1	I	240	PRO	3.3
1	E	867	GLN	3.3
1	I	794	GLN	3.3
1	C	576	ALA	3.3
1	C	241	PHE	3.3
1	G	579	PHE	3.3
1	I	799	PRO	3.3
1	K	575	ARG	3.2
1	E	853	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
1	G	491	LYS	3.2
1	K	329	ALA	3.2
1	C	800	VAL	3.2
1	E	866	ARG	3.2
1	I	859	ARG	3.2
1	C	473	TYR	3.2
1	K	580	ASP	3.1
1	K	577	ALA	3.1
1	E	540	LEU	3.1
1	E	298	ASN	3.1
1	E	397	GLY	3.1
1	K	921	ALA	3.1
1	C	794	GLN	3.0
1	K	862	ILE	3.0
1	E	901	VAL	3.0
1	K	597	LEU	3.0
1	E	607	GLU	3.0
1	K	606	ASN	3.0
1	E	469	ASP	3.0
1	E	481	ASP	3.0
1	K	493	LEU	3.0
1	A	398	SER	3.0
1	E	797	PRO	3.0
1	A	528	LYS	2.9
1	A	502	ARG	2.9
1	G	423	ALA	2.9
1	E	579	PHE	2.9
1	K	607	GLU	2.8
1	I	795	GLU	2.8
1	K	469	ASP	2.8
1	G	717	TYR	2.8
1	I	564	LEU	2.8
1	I	400	GLY	2.8
1	A	800	VAL	2.8
1	C	490	CYS	2.8
1	G	575	ARG	2.8
1	E	796	LYS	2.8
1	A	468	SER	2.8
1	K	717	TYR	2.8
1	K	574	VAL	2.7
1	C	467	ILE	2.7
1	I	526	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	K	473	TYR	2.7
1	E	525	MET	2.7
1	I	468	SER	2.7
1	G	840	GLU	2.7
1	G	489	ILE	2.7
1	A	793	SER	2.7
1	E	272	PHE	2.7
1	I	492	ASP	2.7
1	I	580	ASP	2.7
1	C	564	LEU	2.6
1	E	512	TRP	2.6
1	E	896	ASP	2.6
1	E	601	SER	2.6
1	C	718	VAL	2.6
1	E	868	ASN	2.6
1	E	599	SER	2.6
1	K	477	GLN	2.6
1	E	330	GLU	2.5
1	C	609	PRO	2.5
1	C	789	PHE	2.5
1	E	331	ASN	2.5
1	A	556	GLU	2.5
1	I	509	TYR	2.5
1	C	745	ALA	2.5
1	C	692	ASN	2.5
1	E	879	TYR	2.5
1	A	789	PHE	2.5
1	E	577	ALA	2.5
1	C	556	GLU	2.5
1	K	604	PRO	2.5
1	A	692	ASN	2.4
1	G	467	ILE	2.4
1	A	921	ALA	2.4
1	C	466	ARG	2.4
1	A	484	SER	2.4
1	A	799	PRO	2.4
1	G	298	ASN	2.4
1	E	555	SER	2.4
1	I	398	SER	2.4
1	E	592	GLU	2.4
1	I	595	GLN	2.4
1	G	800	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	575	ARG	2.4
1	C	489	ILE	2.4
1	C	897	ILE	2.4
1	G	416	ASP	2.4
1	C	491	LYS	2.4
1	A	794	GLN	2.4
1	G	272	PHE	2.4
1	G	576	ALA	2.4
1	C	479	MET	2.3
1	E	576	ALA	2.3
1	E	899	THR	2.3
1	E	692	ASN	2.3
1	I	399	SER	2.3
1	I	512	TRP	2.3
1	A	678	GLN	2.3
1	A	491	LYS	2.3
1	C	484	SER	2.3
1	G	601	SER	2.3
1	I	599	SER	2.3
1	K	468	SER	2.3
1	K	481	ASP	2.3
1	E	480	ARG	2.3
1	E	604	PRO	2.3
1	K	609	PRO	2.3
1	G	852	GLN	2.3
1	E	574	VAL	2.3
1	K	680	CYS	2.3
1	E	863	PHE	2.3
1	G	776	PHE	2.3
1	G	242	LYS	2.3
1	G	723	LYS	2.3
1	I	315	GLU	2.3
1	A	488	ARG	2.3
1	E	850	PRO	2.3
1	I	793	SER	2.3
1	E	507	GLN	2.3
1	K	675	LEU	2.3
1	E	895	GLU	2.2
1	K	298	ASN	2.2
1	E	799	PRO	2.2
1	G	398	SER	2.2
1	I	287	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	492	ASP	2.2
1	E	416	ASP	2.2
1	K	472	LYS	2.2
1	G	856	PHE	2.2
1	K	581	GLU	2.2
1	E	564	LEU	2.2
1	K	923	MET	2.2
1	E	877	VAL	2.2
1	E	486	ALA	2.1
1	C	298	ASN	2.1
1	G	794	GLN	2.1
1	I	606	ASN	2.1
1	K	509	TYR	2.1
1	I	288	GLY	2.1
2	B	10	A	2.1
1	C	534	ILE	2.1
1	K	476	ALA	2.1
1	C	548	TYR	2.1
1	G	564	LEU	2.1
1	C	689	GLY	2.1
1	A	580	ASP	2.1
1	E	921	ALA	2.1
1	G	316	ARG	2.1
1	I	502	ARG	2.1
1	C	582	ILE	2.1
1	I	574	VAL	2.1
1	C	580	ASP	2.1
1	E	527	ASP	2.1
1	K	576	ALA	2.1
1	E	513	ILE	2.1
1	I	467	ILE	2.1
1	K	471	PHE	2.1
1	E	795	GLU	2.0
1	E	804	GLU	2.0
1	I	577	ALA	2.0
1	I	592	GLU	2.0
1	K	486	ALA	2.0
1	G	799	PRO	2.0
1	K	678	GLN	2.0
1	A	581	GLU	2.0
1	E	315	GLU	2.0
1	E	590	PHE	2.0

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Mol	Chain	Res	Type	RSRZ
1	E	603	ASP	2.0
1	K	603	ASP	2.0
1	C	855	SER	2.0
1	E	506	THR	2.0
1	E	594	LEU	2.0
1	A	397	GLY	2.0
1	I	692	ASN	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
6	BU3	A	1005	6/6	0.68	0.29	61,75,81,93	0
5	ETF	K	1003	6/6	0.77	0.26	85,120,127,128	0
6	BU3	C	1005	6/6	0.77	0.39	86,100,104,112	0
6	BU3	G	1004	6/6	0.78	0.34	87,98,107,110	0
6	BU3	C	1003	6/6	0.79	0.30	48,70,78,87	0
5	ETF	E	1002	6/6	0.80	0.21	50,106,115,120	0
6	BU3	C	1006	6/6	0.83	0.24	57,65,67,75	0
6	BU3	E	1004	6/6	0.83	0.27	72,85,93,97	0
3	ZN	G	1001	1/1	0.83	0.37	210,210,210,210	0
6	BU3	I	1004	6/6	0.84	0.24	39,61,73,82	0
4	MG	I	1003	1/1	0.85	0.15	57,57,57,57	0
4	MG	I	1002	1/1	0.85	0.16	51,51,51,51	0
6	BU3	G	1005	6/6	0.86	0.23	71,77,88,95	0
6	BU3	G	1003	6/6	0.86	0.24	42,62,76,83	0
6	BU3	K	1004	6/6	0.86	0.20	51,52,61,68	0
4	MG	K	1002	1/1	0.87	0.18	62,62,62,62	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	ETF	G	1002	6/6	0.88	0.26	50,83,90,108	0
6	BU3	A	1004	6/6	0.88	0.26	59,71,77,86	0
6	BU3	E	1003	6/6	0.89	0.21	69,71,75,88	0
6	BU3	C	1004	6/6	0.92	0.24	17,56,70,73	0
5	ETF	A	1003	6/6	0.92	0.19	37,69,80,82	0
4	MG	C	1002	1/1	0.94	0.23	56,56,56,56	0
4	MG	A	1002	1/1	0.96	0.18	44,44,44,44	0
4	MG	H	101	1/1	0.96	0.14	46,46,46,46	0
3	ZN	K	1001	1/1	0.97	0.13	94,94,94,94	0
3	ZN	E	1001	1/1	0.97	0.09	95,95,95,95	0
3	ZN	A	1001	1/1	0.98	0.14	97,97,97,97	0
3	ZN	C	1001	1/1	0.98	0.18	99,99,99,99	0
3	ZN	I	1001	1/1	0.99	0.17	102,102,102,102	0

6.5 Other polymers [i](#)

There are no such residues in this entry.