



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:14 AM UTC

PDB ID : 3P6Y / pdb_00003p6y
Title : CF Im25-CF Im68-UGUAA complex
Authors : Li, H.; Tong, S.; Li, X.; Shi, H.; Gao, Y.; Ge, H.; Niu, L.; Teng, M.
Deposited on : 2010-10-11
Resolution : 2.90 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

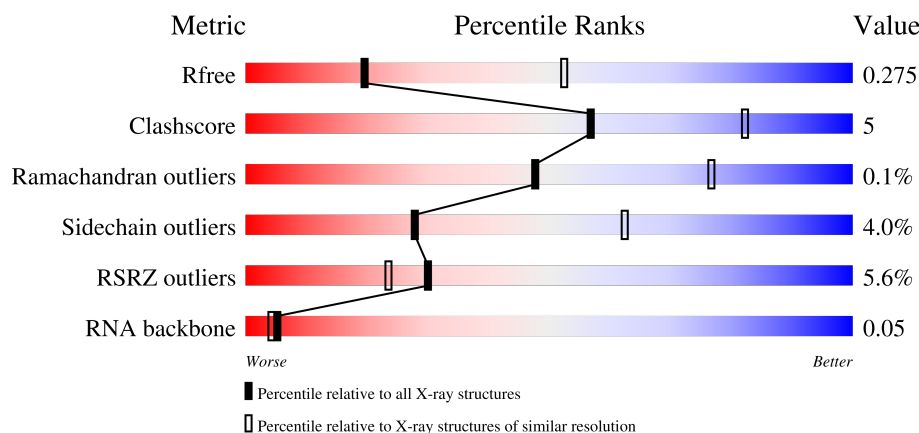
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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.



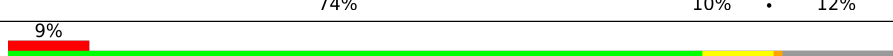
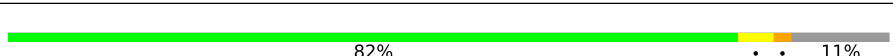
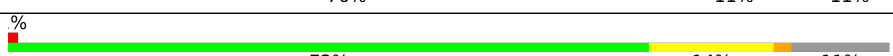
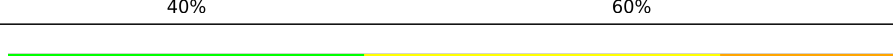
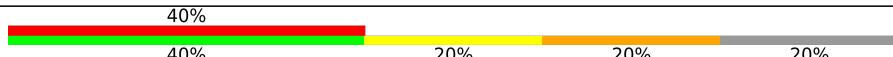
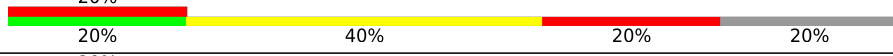
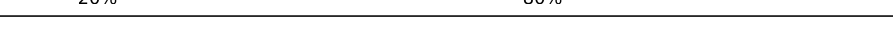
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)
RNA backbone	3983	1120 (3.10-2.70)

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	202	2% 87% 9% .
1	B	202	% 86% 11% ..
1	E	202	2% 85% 11% .
1	F	202	89% 8% .

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Mol	Chain	Length	Quality of chain
1	I	202	
1	J	202	
1	M	202	
1	N	202	
2	C	90	
2	D	90	
2	G	90	
2	H	90	
2	K	90	
2	L	90	
2	O	90	
2	P	90	
3	Q	5	
3	R	5	
3	S	5	
3	T	5	
3	U	5	
3	V	5	
3	W	5	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17567 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 Cleavage and polyadenylation specificity factor subunit 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	0	0
			1568	1022	267	275	4			
1	B	197	Total	C	N	O	S	0	0	0
			1579	1028	268	279	4			
1	E	195	Total	C	N	O	S	0	0	0
			1571	1024	267	276	4			
1	F	197	Total	C	N	O	S	0	0	0
			1583	1033	265	281	4			
1	I	195	Total	C	N	O	S	0	0	0
			1553	1014	266	269	4			
1	J	197	Total	C	N	O	S	0	0	0
			1581	1033	265	279	4			
1	M	196	Total	C	N	O	S	0	0	0
			1570	1023	268	275	4			
1	N	197	Total	C	N	O	S	0	0	0
			1580	1029	268	279	4			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	228	LEU	-	expression tag	UNP O43809
A	229	GLU	-	expression tag	UNP O43809
A	230	HIS	-	expression tag	UNP O43809
A	231	HIS	-	expression tag	UNP O43809
A	232	HIS	-	expression tag	UNP O43809
A	233	HIS	-	expression tag	UNP O43809
A	234	HIS	-	expression tag	UNP O43809
A	235	HIS	-	expression tag	UNP O43809
B	228	LEU	-	expression tag	UNP O43809
B	229	GLU	-	expression tag	UNP O43809
B	230	HIS	-	expression tag	UNP O43809
B	231	HIS	-	expression tag	UNP O43809
B	232	HIS	-	expression tag	UNP O43809

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Chain	Residue	Modelled	Actual	Comment	Reference
B	233	HIS	-	expression tag	UNP O43809
B	234	HIS	-	expression tag	UNP O43809
B	235	HIS	-	expression tag	UNP O43809
E	228	LEU	-	expression tag	UNP O43809
E	229	GLU	-	expression tag	UNP O43809
E	230	HIS	-	expression tag	UNP O43809
E	231	HIS	-	expression tag	UNP O43809
E	232	HIS	-	expression tag	UNP O43809
E	233	HIS	-	expression tag	UNP O43809
E	234	HIS	-	expression tag	UNP O43809
E	235	HIS	-	expression tag	UNP O43809
F	228	LEU	-	expression tag	UNP O43809
F	229	GLU	-	expression tag	UNP O43809
F	230	HIS	-	expression tag	UNP O43809
F	231	HIS	-	expression tag	UNP O43809
F	232	HIS	-	expression tag	UNP O43809
F	233	HIS	-	expression tag	UNP O43809
F	234	HIS	-	expression tag	UNP O43809
F	235	HIS	-	expression tag	UNP O43809
I	228	LEU	-	expression tag	UNP O43809
I	229	GLU	-	expression tag	UNP O43809
I	230	HIS	-	expression tag	UNP O43809
I	231	HIS	-	expression tag	UNP O43809
I	232	HIS	-	expression tag	UNP O43809
I	233	HIS	-	expression tag	UNP O43809
I	234	HIS	-	expression tag	UNP O43809
I	235	HIS	-	expression tag	UNP O43809
J	228	LEU	-	expression tag	UNP O43809
J	229	GLU	-	expression tag	UNP O43809
J	230	HIS	-	expression tag	UNP O43809
J	231	HIS	-	expression tag	UNP O43809
J	232	HIS	-	expression tag	UNP O43809
J	233	HIS	-	expression tag	UNP O43809
J	234	HIS	-	expression tag	UNP O43809
J	235	HIS	-	expression tag	UNP O43809
M	228	LEU	-	expression tag	UNP O43809
M	229	GLU	-	expression tag	UNP O43809
M	230	HIS	-	expression tag	UNP O43809
M	231	HIS	-	expression tag	UNP O43809
M	232	HIS	-	expression tag	UNP O43809
M	233	HIS	-	expression tag	UNP O43809
M	234	HIS	-	expression tag	UNP O43809

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Chain	Residue	Modelled	Actual	Comment	Reference
M	235	HIS	-	expression tag	UNP O43809
N	228	LEU	-	expression tag	UNP O43809
N	229	GLU	-	expression tag	UNP O43809
N	230	HIS	-	expression tag	UNP O43809
N	231	HIS	-	expression tag	UNP O43809
N	232	HIS	-	expression tag	UNP O43809
N	233	HIS	-	expression tag	UNP O43809
N	234	HIS	-	expression tag	UNP O43809
N	235	HIS	-	expression tag	UNP O43809

- Molecule 2 is a protein called Cleavage and polyadenylation specificity factor subunit 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	79	Total	C	N	O	S	0	0	0
			570	366	98	105	1			
2	D	78	Total	C	N	O	S	0	0	0
			557	358	93	105	1			
2	G	80	Total	C	N	O	S	0	0	0
			583	374	101	107	1			
2	H	74	Total	C	N	O		0	0	0
			509	323	88	98				
2	K	80	Total	C	N	O	S	0	0	0
			572	367	99	105	1			
2	L	80	Total	C	N	O	S	0	0	0
			570	367	97	105	1			
2	O	80	Total	C	N	O	S	0	0	0
			581	377	100	103	1			
2	P	75	Total	C	N	O	S	0	0	0
			522	333	89	99	1			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	159	SER	CYS	engineered mutation	UNP Q16630
C	162	LEU	-	expression tag	UNP Q16630
C	163	GLU	-	expression tag	UNP Q16630
C	164	HIS	-	expression tag	UNP Q16630
C	165	HIS	-	expression tag	UNP Q16630
C	166	HIS	-	expression tag	UNP Q16630
C	167	HIS	-	expression tag	UNP Q16630
C	168	HIS	-	expression tag	UNP Q16630
C	169	HIS	-	expression tag	UNP Q16630

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Chain	Residue	Modelled	Actual	Comment	Reference
D	159	SER	CYS	engineered mutation	UNP Q16630
D	162	LEU	-	expression tag	UNP Q16630
D	163	GLU	-	expression tag	UNP Q16630
D	164	HIS	-	expression tag	UNP Q16630
D	165	HIS	-	expression tag	UNP Q16630
D	166	HIS	-	expression tag	UNP Q16630
D	167	HIS	-	expression tag	UNP Q16630
D	168	HIS	-	expression tag	UNP Q16630
D	169	HIS	-	expression tag	UNP Q16630
G	159	SER	CYS	engineered mutation	UNP Q16630
G	162	LEU	-	expression tag	UNP Q16630
G	163	GLU	-	expression tag	UNP Q16630
G	164	HIS	-	expression tag	UNP Q16630
G	165	HIS	-	expression tag	UNP Q16630
G	166	HIS	-	expression tag	UNP Q16630
G	167	HIS	-	expression tag	UNP Q16630
G	168	HIS	-	expression tag	UNP Q16630
G	169	HIS	-	expression tag	UNP Q16630
H	159	SER	CYS	engineered mutation	UNP Q16630
H	162	LEU	-	expression tag	UNP Q16630
H	163	GLU	-	expression tag	UNP Q16630
H	164	HIS	-	expression tag	UNP Q16630
H	165	HIS	-	expression tag	UNP Q16630
H	166	HIS	-	expression tag	UNP Q16630
H	167	HIS	-	expression tag	UNP Q16630
H	168	HIS	-	expression tag	UNP Q16630
H	169	HIS	-	expression tag	UNP Q16630
K	159	SER	CYS	engineered mutation	UNP Q16630
K	162	LEU	-	expression tag	UNP Q16630
K	163	GLU	-	expression tag	UNP Q16630
K	164	HIS	-	expression tag	UNP Q16630
K	165	HIS	-	expression tag	UNP Q16630
K	166	HIS	-	expression tag	UNP Q16630
K	167	HIS	-	expression tag	UNP Q16630
K	168	HIS	-	expression tag	UNP Q16630
K	169	HIS	-	expression tag	UNP Q16630
L	159	SER	CYS	engineered mutation	UNP Q16630
L	162	LEU	-	expression tag	UNP Q16630
L	163	GLU	-	expression tag	UNP Q16630
L	164	HIS	-	expression tag	UNP Q16630
L	165	HIS	-	expression tag	UNP Q16630
L	166	HIS	-	expression tag	UNP Q16630

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Chain	Residue	Modelled	Actual	Comment	Reference
L	167	HIS	-	expression tag	UNP Q16630
L	168	HIS	-	expression tag	UNP Q16630
L	169	HIS	-	expression tag	UNP Q16630
O	159	SER	CYS	engineered mutation	UNP Q16630
O	162	LEU	-	expression tag	UNP Q16630
O	163	GLU	-	expression tag	UNP Q16630
O	164	HIS	-	expression tag	UNP Q16630
O	165	HIS	-	expression tag	UNP Q16630
O	166	HIS	-	expression tag	UNP Q16630
O	167	HIS	-	expression tag	UNP Q16630
O	168	HIS	-	expression tag	UNP Q16630
O	169	HIS	-	expression tag	UNP Q16630
P	159	SER	CYS	engineered mutation	UNP Q16630
P	162	LEU	-	expression tag	UNP Q16630
P	163	GLU	-	expression tag	UNP Q16630
P	164	HIS	-	expression tag	UNP Q16630
P	165	HIS	-	expression tag	UNP Q16630
P	166	HIS	-	expression tag	UNP Q16630
P	167	HIS	-	expression tag	UNP Q16630
P	168	HIS	-	expression tag	UNP Q16630
P	169	HIS	-	expression tag	UNP Q16630

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	Q	2	Total	C	N	O	P	0	0	0
			32	14	7	10	1			
3	R	5	Total	C	N	O	P	0	0	0
			89	42	14	30	3			
3	S	5	Total	C	N	O	P	0	0	0
			104	48	19	33	4			
3	T	4	Total	C	N	O	P	0	0	0
			77	37	14	24	2			
3	U	4	Total	C	N	O	P	0	0	0
			70	33	14	21	2			
3	V	3	Total	C	N	O	P	0	0	0
			60	28	9	21	2			
3	W	1	Total	C	N	O		0	0	0
			17	9	2	6				

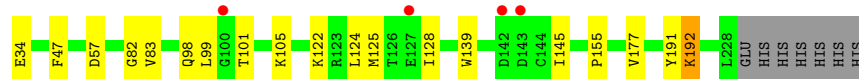
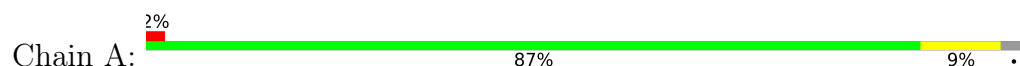
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total O 2 2	0	0
4	B	3	Total O 3 3	0	0
4	C	3	Total O 3 3	0	0
4	D	4	Total O 4 4	0	0
4	E	3	Total O 3 3	0	0
4	F	9	Total O 9 9	0	0
4	G	7	Total O 7 7	0	0
4	H	4	Total O 4 4	0	0
4	I	3	Total O 3 3	0	0
4	J	7	Total O 7 7	0	0
4	K	4	Total O 4 4	0	0
4	L	3	Total O 3 3	0	0
4	M	6	Total O 6 6	0	0
4	N	7	Total O 7 7	0	0
4	O	3	Total O 3 3	0	0
4	P	1	Total O 1 1	0	0

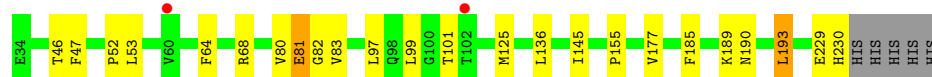
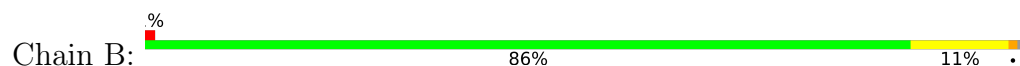
3 Residue-property plots [i](#)

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.

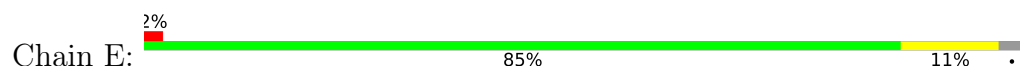
- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5



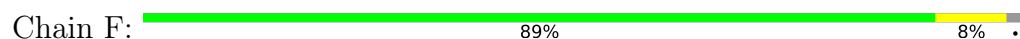
- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5



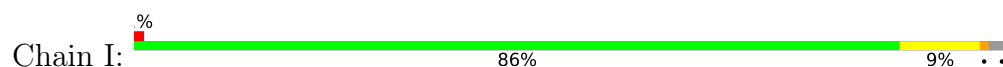
- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5



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- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5

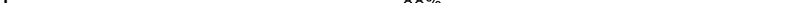


- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5

Chain J: 86% 11%

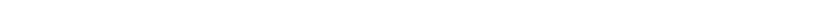


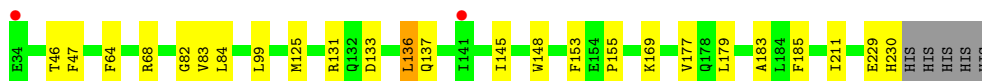
- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5

Chain M:  7% 88% 8% ..

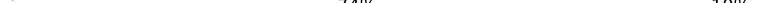


- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5

Chain N:  %

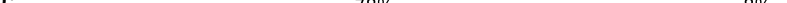


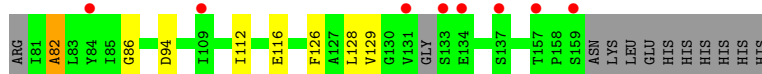
- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6

Chain C:  74% 10% 12%

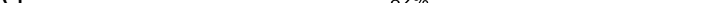


- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6

Chain D:  9% 78% 8% 13%

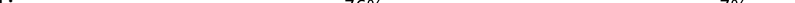


- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6

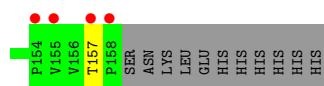
Chain G:  82% . . 11%



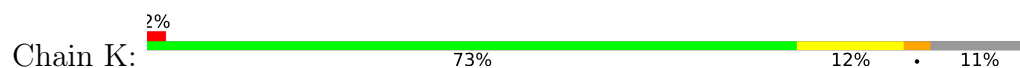
- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6

Chain H: 

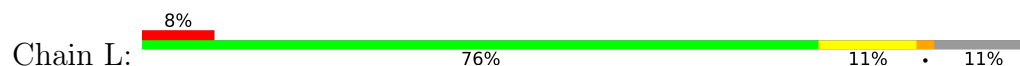




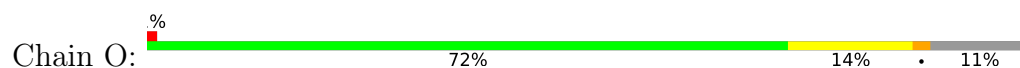
- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6



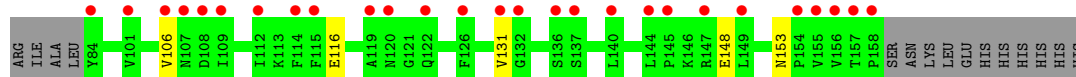
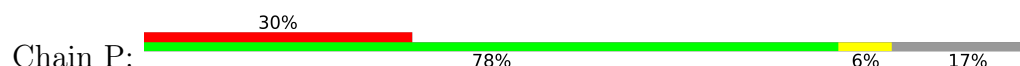
- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6



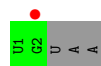
- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6



- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6



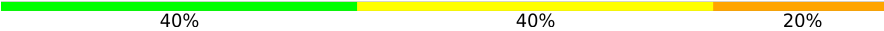
- Molecule 3: 5'-R(*UP*GP*UP*AP*A)-3'



- Molecule 3: 5'-R(*UP*GP*UP*AP*A)-3'

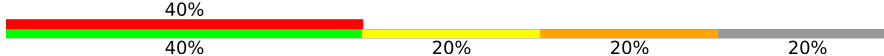


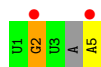
- Molecule 3: 5'-R(*UP*GP*UP*AP*A)-3'

Chain S: 



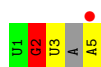
• Molecule 3: 5'-R(*UP*GP*UP*AP*A)-3'

Chain T: 



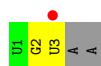
• Molecule 3: 5'-R(*UP*GP*UP*AP*A)-3'

Chain U: 



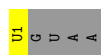
• Molecule 3: 5'-R(*UP*GP*UP*AP*A)-3'

Chain V: 



• Molecule 3: 5'-R(*UP*GP*UP*AP*A)-3'

Chain W: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	104.78Å 129.42Å 111.16Å 90.00° 94.01° 90.00°	Depositor
Resolution (Å)	50.00 – 2.90 50.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.9 (50.00-2.90) 95.9 (50.00-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.12 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.226 , 0.275 0.226 , 0.275	Depositor DCC
R_{free} test set	3182 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	59.3	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	17567	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.65 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.4953e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.44	0/1614	0.76	1/2196 (0.0%)
1	B	0.44	0/1626	0.79	1/2215 (0.0%)
1	E	0.44	0/1617	0.75	2/2200 (0.1%)
1	F	0.44	0/1630	0.75	1/2221 (0.0%)
1	I	0.42	0/1599	0.72	1/2177 (0.0%)
1	J	0.45	0/1628	0.74	1/2218 (0.0%)
1	M	0.44	0/1616	0.74	2/2200 (0.1%)
1	N	0.45	0/1627	0.78	1/2216 (0.0%)
2	C	0.47	0/582	0.76	0/795
2	D	0.46	0/568	0.77	1/776 (0.1%)
2	G	0.43	0/595	0.72	0/812
2	H	0.43	0/519	0.68	0/711
2	K	0.48	0/584	0.80	0/799
2	L	0.43	0/582	0.75	0/796
2	O	0.46	0/593	0.73	0/810
2	P	0.43	0/534	0.70	0/733
3	Q	0.61	0/34	0.58	0/50
3	R	0.46	0/98	1.15	1/149 (0.7%)
3	S	0.46	0/116	0.87	0/179
3	T	0.55	0/85	0.71	0/129
3	U	0.53	0/77	1.17	1/116 (0.9%)
3	V	0.51	0/66	0.85	0/101
3	W	0.63	0/18	0.89	0/26
All	All	0.45	0/18008	0.76	13/24625 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	155	PRO	N-CA-C	6.59	118.74	110.70
1	F	155	PRO	N-CA-C	6.58	118.72	110.70
1	M	155	PRO	N-CA-C	6.41	118.52	110.70
1	B	155	PRO	N-CA-C	6.33	118.42	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	M	101	THR	N-CA-C	-6.15	105.57	114.12
1	E	155	PRO	N-CA-C	5.94	117.95	110.70
1	N	155	PRO	N-CA-C	5.74	117.70	110.70
1	A	155	PRO	N-CA-C	5.46	117.36	110.70
3	R	2	G	C1'-O4'-C4'	-5.40	104.30	109.70
2	D	82	ALA	N-CA-C	5.25	118.04	109.59
3	U	2	G	N9-C1'-C2'	5.24	119.87	112.00
1	E	101	THR	N-CA-C	-5.23	107.92	114.56
1	I	155	PRO	N-CA-C	5.07	116.88	110.70

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	1568	0	1533	11	0
1	B	1579	0	1521	12	0
1	E	1571	0	1534	13	0
1	F	1583	0	1532	13	0
1	I	1553	0	1509	11	0
1	J	1581	0	1532	10	0
1	M	1570	0	1528	16	0
1	N	1580	0	1523	15	0
2	C	570	0	528	8	0
2	D	557	0	493	3	0
2	G	583	0	546	5	0
2	H	509	0	412	2	0
2	K	572	0	523	9	0
2	L	570	0	519	7	0
2	O	581	0	557	9	0
2	P	522	0	448	1	0
3	Q	32	0	15	0	0
3	R	89	0	48	4	0
3	S	104	0	55	3	0
3	T	77	0	42	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	U	70	0	37	5	0
3	V	60	0	33	0	0
3	W	17	0	12	5	0
4	A	2	0	0	0	0
4	B	3	0	0	0	0
4	C	3	0	0	0	0
4	D	4	0	0	0	0
4	E	3	0	0	0	0
4	F	9	0	0	0	0
4	G	7	0	0	0	0
4	H	4	0	0	0	0
4	I	3	0	0	1	0
4	J	7	0	0	2	0
4	K	4	0	0	0	0
4	L	3	0	0	0	0
4	M	6	0	0	1	0
4	N	7	0	0	0	0
4	O	3	0	0	0	0
4	P	1	0	0	0	0
All	All	17567	0	16480	153	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 (153) 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)
3:U:2:G:H1'	3:U:5:A:N6	1.54	1.21
2:K:118:ARG:HH11	2:K:118:ARG:HG3	1.22	1.03
2:G:118:ARG:HH11	2:G:118:ARG:HG3	1.24	1.01
2:C:118:ARG:HH11	2:C:118:ARG:HG3	1.27	0.99
3:U:2:G:H1'	3:U:5:A:H61	1.22	0.92
3:U:2:G:C1'	3:U:5:A:N6	2.33	0.91
2:K:159:SER:HB2	2:K:160:ASN:HA	1.57	0.85
2:O:118:ARG:HH11	2:O:118:ARG:HG3	1.40	0.84
2:O:118:ARG:HH11	2:O:118:ARG:CG	1.95	0.80
2:G:118:ARG:HH11	2:G:118:ARG:CG	1.93	0.79
1:F:104:PHE:O	3:W:1:U:N3	2.18	0.75
1:N:229:GLU:O	1:N:230:HIS:HB2	1.86	0.75
2:K:118:ARG:HH11	2:K:118:ARG:CG	1.97	0.74
1:E:82:GLY:HA3	1:E:125:MET:CE	2.18	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:99:LEU:HD11	1:M:191:TYR:CE2	2.26	0.70
1:B:125:MET:HE3	1:B:177:VAL:HG21	1.73	0.69
3:U:2:G:C1'	3:U:5:A:H61	1.96	0.69
2:L:80:ARG:HB2	3:S:4:A:H61	1.57	0.69
1:A:82:GLY:HA3	1:A:125:MET:CE	2.23	0.69
1:E:124:LEU:O	1:E:128:ILE:HG12	1.94	0.68
3:T:2:G:H21	3:T:5:A:H62	1.43	0.67
2:C:118:ARG:HG3	2:C:118:ARG:NH1	2.02	0.65
2:G:118:ARG:HG3	2:G:118:ARG:NH1	2.04	0.65
1:B:83:VAL:HG21	1:B:145:ILE:HD11	1.79	0.64
1:N:64:PHE:O	1:N:68:ARG:HG3	1.98	0.63
1:M:82:GLY:HA3	1:M:125:MET:CE	2.29	0.62
2:D:82:ALA:HB1	2:D:128:LEU:HD11	1.81	0.62
1:J:142:ASP:HB3	4:J:33:HOH:O	2.00	0.62
1:N:82:GLY:HA3	1:N:125:MET:CE	2.29	0.62
1:F:82:GLY:HA3	1:F:125:MET:CE	2.29	0.62
1:I:82:GLY:HA3	1:I:125:MET:CE	2.29	0.62
2:O:159:SER:O	2:O:160:ASN:HB2	2.00	0.61
1:F:104:PHE:O	3:W:1:U:C2	2.54	0.61
1:B:229:GLU:O	1:B:230:HIS:HB2	2.02	0.60
3:R:3:U:H5''	3:R:3:U:H6	1.66	0.60
1:N:137:GLN:HE22	1:N:183:ALA:HB1	1.67	0.60
1:E:82:GLY:HA3	1:E:125:MET:HE1	1.82	0.60
1:F:125:MET:HE3	1:F:177:VAL:HG21	1.84	0.58
1:A:122:LYS:HG2	1:A:139:TRP:HB2	1.86	0.58
1:M:142:ASP:HB3	4:M:32:HOH:O	2.03	0.57
3:R:4:A:O3'	3:R:5:A:H3'	2.03	0.57
1:F:64:PHE:O	1:F:68:ARG:HG3	2.05	0.56
1:B:52:PRO:HG3	1:B:189:LYS:HB2	1.86	0.56
1:M:118:VAL:HG13	1:M:141:ILE:HG12	1.88	0.55
1:I:82:GLY:HA3	1:I:125:MET:HE1	1.88	0.54
1:N:125:MET:HE3	1:N:177:VAL:HG21	1.90	0.54
2:K:118:ARG:HG3	2:K:118:ARG:NH1	2.02	0.54
2:O:98:THR:HG22	2:O:102:HIS:CE1	2.43	0.54
1:A:98:GLN:NE2	1:A:101:THR:HA	2.23	0.54
1:A:98:GLN:HE21	1:A:101:THR:HA	1.73	0.54
2:G:153:ASN:HD22	2:G:153:ASN:N	2.06	0.54
1:A:47:PHE:O	1:N:133:ASP:HB2	2.07	0.53
2:O:159:SER:O	2:O:160:ASN:CB	2.56	0.53
1:M:97:LEU:HB3	1:M:191:TYR:HD2	1.73	0.53
1:B:47:PHE:CE1	1:B:185:PHE:HD2	2.26	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:118:VAL:HG22	1:M:141:ILE:HD11	1.92	0.52
3:R:1:U:H6	3:R:1:U:H5''	1.75	0.52
1:J:64:PHE:O	1:J:68:ARG:HG3	2.10	0.52
1:A:124:LEU:O	1:A:128:ILE:HG12	2.11	0.51
1:E:95:LEU:HD22	1:E:193:LEU:HD21	1.93	0.51
1:M:59:SER:OG	1:M:62:ALA:N	2.44	0.51
1:F:131:ARG:HD3	1:F:133:ASP:OD1	2.12	0.50
1:A:105:LYS:HE3	1:A:191:TYR:HE2	1.76	0.50
1:A:192:LYS:HB3	1:N:136:LEU:HB3	1.92	0.50
2:L:101:VAL:HG11	2:L:109:ILE:HD11	1.93	0.50
2:C:112:ILE:HG12	2:C:129:VAL:HG22	1.93	0.50
1:J:83:VAL:HG21	1:J:145:ILE:HD11	1.93	0.50
1:B:53:LEU:O	1:B:190:ASN:ND2	2.42	0.49
1:N:82:GLY:HA3	1:N:125:MET:HE1	1.93	0.49
1:M:46:THR:HB	1:M:184:LEU:HD12	1.95	0.49
1:B:82:GLY:HA3	1:B:125:MET:CE	2.41	0.49
1:E:125:MET:HE3	1:E:177:VAL:HG21	1.95	0.49
2:O:118:ARG:CG	2:O:118:ARG:NH1	2.63	0.49
1:B:82:GLY:HA3	1:B:125:MET:HE1	1.92	0.49
2:L:139:LYS:O	2:L:143:LEU:HG	2.11	0.49
2:L:80:ARG:CB	3:S:4:A:H61	2.26	0.49
1:N:47:PHE:CE1	1:N:185:PHE:HD2	2.31	0.49
1:M:99:LEU:CD1	1:M:191:TYR:CE2	2.95	0.48
1:A:125:MET:HE3	1:A:177:VAL:HG21	1.94	0.48
2:C:118:ARG:NH1	2:C:118:ARG:CG	2.74	0.48
1:N:131:ARG:HD3	1:N:133:ASP:OD1	2.14	0.48
2:K:118:ARG:CG	2:K:118:ARG:NH1	2.64	0.48
1:F:83:VAL:HG21	1:F:145:ILE:HD11	1.95	0.48
1:I:105:LYS:HE3	1:I:191:TYR:CE2	2.49	0.48
3:U:2:G:C1'	3:U:5:A:H62	2.23	0.47
1:I:124:LEU:O	1:I:128:ILE:HG12	2.13	0.47
1:B:64:PHE:O	1:B:68:ARG:HG3	2.15	0.47
1:M:82:GLY:HA3	1:M:125:MET:HE3	1.96	0.47
1:E:99:LEU:HD13	1:E:191:TYR:CE2	2.49	0.47
1:F:104:PHE:O	3:W:1:U:O2	2.32	0.47
1:J:121:LEU:HD11	1:J:125:MET:HE2	1.96	0.47
2:L:80:ARG:HB2	3:S:4:A:N6	2.28	0.47
1:B:80:VAL:C	1:B:81:GLU:HG2	2.39	0.47
1:M:97:LEU:HD11	1:M:128:ILE:HG21	1.97	0.47
1:I:125:MET:HE3	1:I:177:VAL:HG21	1.97	0.46
1:F:104:PHE:N	3:W:1:U:O2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:135:ALA:O	2:H:139:LYS:CB	2.63	0.46
1:J:97:LEU:HD21	1:J:129:LEU:HD21	1.98	0.46
3:T:2:G:N2	3:T:5:A:H62	2.10	0.46
1:A:83:VAL:HG21	1:A:145:ILE:HD11	1.97	0.46
1:N:83:VAL:HG21	1:N:145:ILE:HD11	1.98	0.46
2:H:84:TYR:N	2:H:157:THR:O	2.49	0.46
2:K:96:ASP:OD2	2:K:150:HIS:HE1	1.99	0.45
1:M:59:SER:OG	1:M:62:ALA:CB	2.64	0.45
1:J:142:ASP:HB2	4:J:27:HOH:O	2.16	0.45
2:C:83:LEU:HD23	2:C:140:LEU:HD23	1.99	0.45
1:J:82:GLY:HA3	1:J:125:MET:CE	2.47	0.45
2:L:112:ILE:HG12	2:L:129:VAL:HG22	1.99	0.45
1:I:81:GLU:HG2	1:I:106:LEU:HB2	1.99	0.44
1:I:188:PRO:HB2	1:I:191:TYR:HD1	1.81	0.44
1:N:137:GLN:NE2	1:N:183:ALA:HB1	2.32	0.44
1:I:95:LEU:HD22	1:I:193:LEU:HD21	2.00	0.44
1:I:215:LEU:N	1:I:216:PRO:CD	2.81	0.44
1:I:52:PRO:HG3	1:I:189:LYS:HB2	2.01	0.43
1:I:142:ASP:HB3	4:I:236:HOH:O	2.19	0.43
1:E:93:HIS:HD2	1:E:195:ALA:O	2.02	0.43
2:D:86:GLY:HA2	2:D:126:PHE:HB3	2.01	0.42
2:O:101:VAL:HG12	2:O:106:VAL:HB	2.00	0.42
2:P:148:GLU:HB3	2:P:153:ASN:ND2	2.34	0.42
2:D:112:ILE:HG12	2:D:129:VAL:HG22	2.00	0.42
1:J:118:VAL:HG13	1:J:141:ILE:HG12	2.00	0.42
1:M:99:LEU:HD11	1:M:191:TYR:CZ	2.54	0.42
2:L:137:SER:O	2:L:141:MET:HG3	2.19	0.42
2:C:97:LEU:O	2:C:101:VAL:HG23	2.19	0.42
1:N:84:LEU:HD23	1:N:179:LEU:HD21	2.01	0.42
1:A:82:GLY:HA3	1:A:125:MET:HE1	2.00	0.42
1:B:47:PHE:CE1	1:B:185:PHE:CD2	3.06	0.42
2:K:153:ASN:HA	2:K:154:PRO:HD3	1.93	0.41
1:J:150:ARG:HD3	1:J:154:GLU:O	2.20	0.41
1:F:82:GLY:HA3	1:F:125:MET:HE1	1.99	0.41
3:R:3:U:H5''	3:R:3:U:C6	2.52	0.41
1:B:97:LEU:HD23	1:B:193:LEU:HD12	2.02	0.41
1:N:153:PHE:CZ	1:N:169:LYS:HD2	2.55	0.41
1:F:102:THR:O	3:W:1:U:O2'	2.30	0.41
1:M:52:PRO:HG3	1:M:189:LYS:HB2	2.03	0.41
1:M:99:LEU:HD11	1:M:191:TYR:CD2	2.54	0.41
2:K:142:ASP:O	2:K:146:LYS:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:91:LEU:HA	1:E:92:PRO:HD3	1.97	0.41
1:F:196:ALA:HA	1:F:197:PRO:HD2	1.94	0.41
1:N:148:TRP:CD2	1:N:211:ILE:HG12	2.56	0.41
2:C:148:GLU:HA	2:C:153:ASN:HD22	1.86	0.41
1:E:51:GLU:HA	1:E:52:PRO:HD3	1.96	0.41
1:E:214:SER:HB2	1:F:217:GLN:HG2	2.03	0.41
2:K:148:GLU:HA	2:K:153:ASN:ND2	2.35	0.41
1:M:99:LEU:CD1	1:M:191:TYR:CD2	3.04	0.41
2:C:138:LYS:HE2	2:C:138:LYS:HB3	1.89	0.41
1:E:105:LYS:HE3	1:E:191:TYR:HE2	1.86	0.41
2:O:144:LEU:HB3	2:O:145:PRO:HD3	2.03	0.41
1:E:47:PHE:CE1	1:E:193:LEU:HD22	2.57	0.40
2:G:157:THR:HA	2:G:158:PRO:HD3	1.93	0.40
1:E:96:LEU:HB2	1:E:194:VAL:HG23	2.03	0.40
2:O:153:ASN:HA	2:O:154:PRO:HD3	1.93	0.40
1:J:121:LEU:HD23	1:J:141:ILE:HD13	2.02	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	193/202 (96%)	186 (96%)	7 (4%)	0	100	100
1	B	195/202 (96%)	190 (97%)	5 (3%)	0	100	100
1	E	193/202 (96%)	188 (97%)	5 (3%)	0	100	100
1	F	195/202 (96%)	190 (97%)	5 (3%)	0	100	100
1	I	193/202 (96%)	186 (96%)	7 (4%)	0	100	100
1	J	195/202 (96%)	189 (97%)	6 (3%)	0	100	100
1	M	194/202 (96%)	184 (95%)	10 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	N	195/202 (96%)	189 (97%)	6 (3%)	0	100	100
2	C	77/90 (86%)	75 (97%)	2 (3%)	0	100	100
2	D	74/90 (82%)	70 (95%)	4 (5%)	0	100	100
2	G	78/90 (87%)	75 (96%)	2 (3%)	1 (1%)	9	32
2	H	68/90 (76%)	67 (98%)	1 (2%)	0	100	100
2	K	78/90 (87%)	75 (96%)	2 (3%)	1 (1%)	9	32
2	L	78/90 (87%)	76 (97%)	2 (3%)	0	100	100
2	O	78/90 (87%)	78 (100%)	0	0	100	100
2	P	73/90 (81%)	70 (96%)	2 (3%)	1 (1%)	9	30
All	All	2157/2336 (92%)	2088 (97%)	66 (3%)	3 (0%)	48	77

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	G	158	PRO
2	K	158	PRO
2	P	106	VAL

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	163/181 (90%)	159 (98%)	4 (2%)	42	74
1	B	162/181 (90%)	156 (96%)	6 (4%)	30	64
1	E	163/181 (90%)	160 (98%)	3 (2%)	51	80
1	F	164/181 (91%)	162 (99%)	2 (1%)	63	86
1	I	158/181 (87%)	153 (97%)	5 (3%)	34	68
1	J	163/181 (90%)	153 (94%)	10 (6%)	17	46
1	M	162/181 (90%)	158 (98%)	4 (2%)	42	74
1	N	162/181 (90%)	159 (98%)	3 (2%)	50	79

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	53/78 (68%)	48 (91%)	5 (9%)	8	26
2	D	49/78 (63%)	47 (96%)	2 (4%)	27	61
2	G	55/78 (70%)	52 (94%)	3 (6%)	19	50
2	H	41/78 (53%)	39 (95%)	2 (5%)	22	54
2	K	52/78 (67%)	46 (88%)	6 (12%)	5	18
2	L	51/78 (65%)	46 (90%)	5 (10%)	7	25
2	O	55/78 (70%)	49 (89%)	6 (11%)	6	20
2	P	45/78 (58%)	43 (96%)	2 (4%)	25	59
All	All	1698/2072 (82%)	1630 (96%)	68 (4%)	28	62

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	GLU
1	A	57	ASP
1	A	99	LEU
1	A	192	LYS
1	B	46	THR
1	B	81	GLU
1	B	99	LEU
1	B	101	THR
1	B	136	LEU
1	B	193	LEU
2	C	83	LEU
2	C	94	ASP
2	C	116	GLU
2	C	118	ARG
2	C	148	GLU
2	D	94	ASP
2	D	116	GLU
1	E	57	ASP
1	E	59	SER
1	E	193	LEU
1	F	99	LEU
1	F	136	LEU
2	G	118	ARG
2	G	148	GLU
2	G	149	LEU
2	H	94	ASP

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Mol	Chain	Res	Type
2	H	116	GLU
1	I	57	ASP
1	I	59	SER
1	I	81	GLU
1	I	110	GLU
1	I	193	LEU
1	J	46	THR
1	J	49	THR
1	J	81	GLU
1	J	99	LEU
1	J	101	THR
1	J	136	LEU
1	J	141	ILE
1	J	193	LEU
1	J	225	ILE
1	J	230	HIS
2	K	116	GLU
2	K	118	ARG
2	K	140	LEU
2	K	148	GLU
2	K	149	LEU
2	K	160	ASN
2	L	80	ARG
2	L	83	LEU
2	L	116	GLU
2	L	131	VAL
2	L	143	LEU
1	M	57	ASP
1	M	101	THR
1	M	119	GLU
1	M	141	ILE
1	N	46	THR
1	N	99	LEU
1	N	136	LEU
2	O	83	LEU
2	O	98	THR
2	O	110	LEU
2	O	116	GLU
2	O	118	ARG
2	O	148	GLU
2	P	116	GLU
2	P	131	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	147	ASN
1	A	152	ASN
1	B	112	ASN
1	B	152	ASN
1	B	157	GLN
1	B	178	GLN
2	C	152	GLN
2	C	153	ASN
2	D	152	GLN
2	D	153	ASN
1	E	65	GLN
1	E	89	HIS
1	E	147	ASN
1	F	157	GLN
2	G	102	HIS
2	G	153	ASN
1	I	89	HIS
1	I	98	GLN
1	I	147	ASN
1	I	152	ASN
1	J	112	ASN
1	J	147	ASN
1	J	178	GLN
2	K	150	HIS
2	K	153	ASN
2	L	153	ASN
1	M	89	HIS
1	M	147	ASN
1	M	152	ASN
1	M	157	GLN
1	N	89	HIS
1	N	112	ASN
2	O	102	HIS
2	O	153	ASN
2	P	153	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	Q	0/5	-	-

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	R	3/5 (60%)	3 (100%)	1 (33%)
3	S	4/5 (80%)	3 (75%)	0
3	T	2/5 (40%)	1 (50%)	0
3	U	2/5 (40%)	2 (100%)	1 (50%)
3	V	2/5 (40%)	1 (50%)	1 (50%)
3	W	0/5	-	-
All	All	13/35 (37%)	10 (76%)	3 (23%)

All (10) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	R	2	G
3	R	3	U
3	R	4	A
3	S	2	G
3	S	4	A
3	S	5	A
3	T	2	G
3	U	2	G
3	U	3	U
3	V	3	U

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	R	3	U
3	U	2	G
3	V	2	G

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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](#)

There are no ligands in this entry.

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	195/202 (96%)	0.16	4 (2%) 63 54	30, 51, 79, 97	0
1	B	197/202 (97%)	-0.02	2 (1%) 79 73	25, 45, 71, 77	0
1	E	195/202 (96%)	0.35	5 (2%) 57 48	34, 57, 81, 97	0
1	F	197/202 (97%)	0.09	0 100 100	29, 52, 82, 89	0
1	I	195/202 (96%)	0.31	3 (1%) 72 64	34, 59, 87, 109	0
1	J	197/202 (97%)	0.11	1 (0%) 87 83	31, 50, 88, 97	0
1	M	196/202 (97%)	0.71	14 (7%) 22 17	35, 65, 104, 117	0
1	N	197/202 (97%)	0.08	2 (1%) 79 73	26, 54, 72, 83	0
2	C	79/90 (87%)	0.03	1 (1%) 75 67	33, 47, 60, 66	0
2	D	78/90 (86%)	0.91	8 (10%) 12 10	50, 82, 95, 115	0
2	G	80/90 (88%)	0.36	0 100 100	29, 56, 78, 86	0
2	H	74/90 (82%)	2.27	41 (55%) 0 0	97, 142, 190, 200	0
2	K	80/90 (88%)	0.27	2 (2%) 58 48	37, 55, 74, 80	0
2	L	80/90 (88%)	0.98	7 (8%) 15 13	53, 79, 105, 117	0
2	O	80/90 (88%)	-0.07	1 (1%) 75 67	29, 46, 61, 80	0
2	P	75/90 (83%)	1.79	27 (36%) 1 0	90, 137, 166, 175	0
3	Q	2/5 (40%)	2.24	1 (50%) 0 0	36, 36, 36, 37	2 (100%)
3	R	5/5 (100%)	1.75	2 (40%) 1 0	63, 73, 76, 83	2 (40%)
3	S	5/5 (100%)	1.46	0 100 100	91, 96, 120, 121	0
3	T	4/5 (80%)	2.28	2 (50%) 0 0	30, 32, 38, 43	4 (100%)
3	U	4/5 (80%)	1.72	1 (25%) 2 1	26, 38, 38, 49	3 (75%)
3	V	3/5 (60%)	1.83	1 (33%) 1 1	25, 25, 31, 37	3 (100%)
3	W	1/5 (20%)	1.93	0 100 100	40, 40, 40, 40	1 (100%)
All	All	2219/2371 (93%)	0.40	125 (5%) 30 23	25, 56, 117, 200	15 (0%)

All (125) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	158	PRO	8.5
2	P	131	VAL	4.7
2	P	158	PRO	4.6
2	H	157	THR	4.5
3	T	5	A	4.5
2	H	133	SER	4.4
2	H	135	ALA	4.3
1	M	133	ASP	4.2
2	H	122	GLN	4.1
2	D	133	SER	4.1
3	R	5	A	4.0
3	V	3	U	4.0
2	P	147	ARG	4.0
1	I	133	ASP	3.8
1	E	143	ASP	3.7
1	M	144	CYS	3.7
2	P	84	TYR	3.7
3	U	5	A	3.7
2	H	84	TYR	3.6
2	P	144	LEU	3.5
2	H	98	THR	3.5
2	H	147	ARG	3.5
2	P	136	SER	3.4
2	P	137	SER	3.4
2	D	137	SER	3.3
3	Q	2	G	3.3
2	H	85	ILE	3.3
1	M	51	GLU	3.2
2	H	149	LEU	3.2
2	H	134	GLU	3.2
2	H	137	SER	3.2
2	H	141	MET	3.2
2	P	119	ALA	3.1
2	P	132	GLY	3.1
2	H	99	GLU	3.1
2	P	149	LEU	3.1
3	R	4	A	3.1
2	H	97	LEU	3.1
2	P	157	THR	3.1
2	H	123	SER	3.1
2	P	154	PRO	3.0
1	E	102	THR	2.9

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Mol	Chain	Res	Type	RSRZ
1	M	187	VAL	2.9
2	P	114	PHE	2.9
2	D	84	TYR	2.9
2	P	108	ASP	2.9
2	H	106	VAL	2.9
2	H	146	LYS	2.8
2	H	119	ALA	2.8
2	H	131	VAL	2.8
2	H	144	LEU	2.8
2	P	106	VAL	2.8
1	M	140	VAL	2.8
2	H	155	VAL	2.8
2	H	154	PRO	2.7
1	E	142	ASP	2.7
1	I	141	ILE	2.7
1	A	143	ASP	2.7
2	H	114	PHE	2.7
2	H	121	GLY	2.7
2	H	115	PHE	2.6
2	H	117	ASN	2.6
1	M	101	THR	2.6
2	P	126	PHE	2.6
1	M	190	ASN	2.6
1	M	50	LYS	2.5
2	H	93	THR	2.5
2	H	120	ASN	2.5
2	D	109	ILE	2.5
2	L	141	MET	2.5
2	H	150	HIS	2.5
2	H	136	SER	2.5
1	B	102	THR	2.5
2	P	101	VAL	2.4
2	P	140	LEU	2.4
2	H	108	ASP	2.4
2	P	120	ASN	2.4
2	H	126	PHE	2.4
2	H	90	TRP	2.4
1	M	53	LEU	2.4
1	N	141	ILE	2.4
2	L	107	ASN	2.3
2	K	136	SER	2.3
1	A	100	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
2	H	83	LEU	2.3
1	N	34	GLU	2.3
1	E	141	ILE	2.3
2	D	157	THR	2.3
2	H	101	VAL	2.3
2	P	155	VAL	2.3
2	L	151	GLY	2.3
2	H	127	ALA	2.3
2	P	156	VAL	2.3
1	I	81	GLU	2.2
1	M	189	LYS	2.2
2	C	108	ASP	2.2
2	L	142	ASP	2.2
1	M	52	PRO	2.2
2	L	155	VAL	2.2
2	D	159	SER	2.2
1	J	99	LEU	2.2
2	H	95	GLU	2.2
2	P	112	ILE	2.2
2	P	122	GLN	2.2
2	P	145	PRO	2.2
2	L	159	SER	2.2
2	P	115	PHE	2.2
2	H	100	ALA	2.2
2	P	109	ILE	2.1
1	M	48	GLY	2.1
2	O	137	SER	2.1
1	B	60	VAL	2.1
1	M	60	VAL	2.1
2	D	131	VAL	2.1
1	A	142	ASP	2.1
1	M	75	GLY	2.1
2	K	134	GLU	2.1
1	E	34	GLU	2.1
2	H	124	LYS	2.1
2	P	107	ASN	2.0
3	T	2	G	2.0
2	H	109	ILE	2.0
1	A	127	GLU	2.0
2	D	134	GLU	2.0
2	L	84	TYR	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.