



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 09:05 AM EDT

PDB ID : 3MX4 / pdb_00003mx4
Title : DNA binding and cleavage by the GIY-YIG endonuclease R.Eco29KI inactive variant E142Q
Authors : Mak, A.N.S.; Lambert, A.R.; Stoddard, B.L.
Deposited on : 2010-05-06
Resolution : 2.50 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
Xtrriage (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

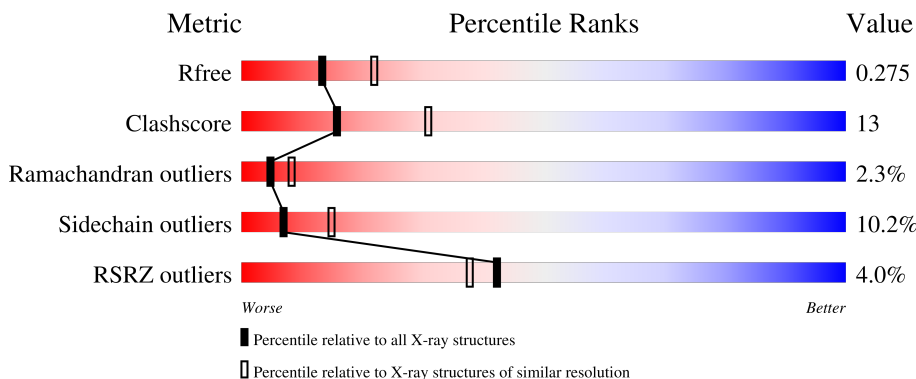
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	235	 4% 66% 20% 9%
1	B	235	 5% 63% 22% 5% 9%
1	C	235	 3% 58% 24% 6% 11%
1	D	235	 3% 60% 22% 6% 11%
1	E	235	 4% 63% 22% 6% 10%

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Mol	Chain	Length	Quality of chain
1	F	235	<p>3% 60% 21% 7% • 11%</p>
1	G	235	<p>2% 63% 19% 5% • 11%</p>
1	H	235	<p>5% 63% 20% 6% 11%</p>
2	I	22	<p>14% 55% 36% 9%</p>
2	K	22	<p>55% 41% 5%</p>
2	M	22	<p>5% 23% 64% 14%</p>
2	O	22	<p>23% 59% 18%</p>
3	J	22	<p>5% 50% 45% 5%</p>
3	L	22	<p>14% 36% 41% 23%</p>
3	N	22	<p>9% 27% 59% 14%</p>
3	P	22	<p>32% 55% 14%</p>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 17406 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 Eco29kIR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	213	Total 1677	C 1069	N 298	O 306	S 4	0	0	0
1	B	213	Total 1687	C 1077	N 299	O 307	S 4	0	0	0
1	C	208	Total 1649	C 1051	N 292	O 302	S 4	0	0	0
1	D	208	Total 1634	C 1041	N 290	O 299	S 4	0	0	0
1	E	211	Total 1663	C 1062	N 297	O 300	S 4	0	0	0
1	F	210	Total 1674	C 1066	N 298	O 306	S 4	0	0	0
1	G	208	Total 1663	C 1062	N 294	O 303	S 4	0	0	0
1	H	209	Total 1658	C 1056	N 295	O 303	S 4	0	0	0

There are 192 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	expression tag	UNP Q46944
A	-19	ALA	-	expression tag	UNP Q46944
A	-18	SER	-	expression tag	UNP Q46944
A	-17	SER	-	expression tag	UNP Q46944
A	-16	HIS	-	expression tag	UNP Q46944
A	-15	HIS	-	expression tag	UNP Q46944
A	-14	HIS	-	expression tag	UNP Q46944
A	-13	HIS	-	expression tag	UNP Q46944
A	-12	HIS	-	expression tag	UNP Q46944
A	-11	HIS	-	expression tag	UNP Q46944
A	-10	SER	-	expression tag	UNP Q46944
A	-9	SER	-	expression tag	UNP Q46944
A	-8	GLY	-	expression tag	UNP Q46944

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	LEU	-	expression tag	UNP Q46944
A	-6	VAL	-	expression tag	UNP Q46944
A	-5	PRO	-	expression tag	UNP Q46944
A	-4	ARG	-	expression tag	UNP Q46944
A	-3	GLY	-	expression tag	UNP Q46944
A	-2	SER	-	expression tag	UNP Q46944
A	-1	SER	-	expression tag	UNP Q46944
A	0	MET	-	expression tag	UNP Q46944
A	1	GLY	-	expression tag	UNP Q46944
A	69	LYS	LEU	engineered mutation	UNP Q46944
A	142	GLN	GLU	engineered mutation	UNP Q46944
B	-20	MET	-	expression tag	UNP Q46944
B	-19	ALA	-	expression tag	UNP Q46944
B	-18	SER	-	expression tag	UNP Q46944
B	-17	SER	-	expression tag	UNP Q46944
B	-16	HIS	-	expression tag	UNP Q46944
B	-15	HIS	-	expression tag	UNP Q46944
B	-14	HIS	-	expression tag	UNP Q46944
B	-13	HIS	-	expression tag	UNP Q46944
B	-12	HIS	-	expression tag	UNP Q46944
B	-11	HIS	-	expression tag	UNP Q46944
B	-10	SER	-	expression tag	UNP Q46944
B	-9	SER	-	expression tag	UNP Q46944
B	-8	GLY	-	expression tag	UNP Q46944
B	-7	LEU	-	expression tag	UNP Q46944
B	-6	VAL	-	expression tag	UNP Q46944
B	-5	PRO	-	expression tag	UNP Q46944
B	-4	ARG	-	expression tag	UNP Q46944
B	-3	GLY	-	expression tag	UNP Q46944
B	-2	SER	-	expression tag	UNP Q46944
B	-1	SER	-	expression tag	UNP Q46944
B	0	MET	-	expression tag	UNP Q46944
B	1	GLY	-	expression tag	UNP Q46944
B	69	LYS	LEU	engineered mutation	UNP Q46944
B	142	GLN	GLU	engineered mutation	UNP Q46944
C	-20	MET	-	expression tag	UNP Q46944
C	-19	ALA	-	expression tag	UNP Q46944
C	-18	SER	-	expression tag	UNP Q46944
C	-17	SER	-	expression tag	UNP Q46944
C	-16	HIS	-	expression tag	UNP Q46944
C	-15	HIS	-	expression tag	UNP Q46944
C	-14	HIS	-	expression tag	UNP Q46944

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-13	HIS	-	expression tag	UNP Q46944
C	-12	HIS	-	expression tag	UNP Q46944
C	-11	HIS	-	expression tag	UNP Q46944
C	-10	SER	-	expression tag	UNP Q46944
C	-9	SER	-	expression tag	UNP Q46944
C	-8	GLY	-	expression tag	UNP Q46944
C	-7	LEU	-	expression tag	UNP Q46944
C	-6	VAL	-	expression tag	UNP Q46944
C	-5	PRO	-	expression tag	UNP Q46944
C	-4	ARG	-	expression tag	UNP Q46944
C	-3	GLY	-	expression tag	UNP Q46944
C	-2	SER	-	expression tag	UNP Q46944
C	-1	SER	-	expression tag	UNP Q46944
C	0	MET	-	expression tag	UNP Q46944
C	1	GLY	-	expression tag	UNP Q46944
C	69	LYS	LEU	engineered mutation	UNP Q46944
C	142	GLN	GLU	engineered mutation	UNP Q46944
D	-20	MET	-	expression tag	UNP Q46944
D	-19	ALA	-	expression tag	UNP Q46944
D	-18	SER	-	expression tag	UNP Q46944
D	-17	SER	-	expression tag	UNP Q46944
D	-16	HIS	-	expression tag	UNP Q46944
D	-15	HIS	-	expression tag	UNP Q46944
D	-14	HIS	-	expression tag	UNP Q46944
D	-13	HIS	-	expression tag	UNP Q46944
D	-12	HIS	-	expression tag	UNP Q46944
D	-11	HIS	-	expression tag	UNP Q46944
D	-10	SER	-	expression tag	UNP Q46944
D	-9	SER	-	expression tag	UNP Q46944
D	-8	GLY	-	expression tag	UNP Q46944
D	-7	LEU	-	expression tag	UNP Q46944
D	-6	VAL	-	expression tag	UNP Q46944
D	-5	PRO	-	expression tag	UNP Q46944
D	-4	ARG	-	expression tag	UNP Q46944
D	-3	GLY	-	expression tag	UNP Q46944
D	-2	SER	-	expression tag	UNP Q46944
D	-1	SER	-	expression tag	UNP Q46944
D	0	MET	-	expression tag	UNP Q46944
D	1	GLY	-	expression tag	UNP Q46944
D	69	LYS	LEU	engineered mutation	UNP Q46944
D	142	GLN	GLU	engineered mutation	UNP Q46944
E	-20	MET	-	expression tag	UNP Q46944

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-19	ALA	-	expression tag	UNP Q46944
E	-18	SER	-	expression tag	UNP Q46944
E	-17	SER	-	expression tag	UNP Q46944
E	-16	HIS	-	expression tag	UNP Q46944
E	-15	HIS	-	expression tag	UNP Q46944
E	-14	HIS	-	expression tag	UNP Q46944
E	-13	HIS	-	expression tag	UNP Q46944
E	-12	HIS	-	expression tag	UNP Q46944
E	-11	HIS	-	expression tag	UNP Q46944
E	-10	SER	-	expression tag	UNP Q46944
E	-9	SER	-	expression tag	UNP Q46944
E	-8	GLY	-	expression tag	UNP Q46944
E	-7	LEU	-	expression tag	UNP Q46944
E	-6	VAL	-	expression tag	UNP Q46944
E	-5	PRO	-	expression tag	UNP Q46944
E	-4	ARG	-	expression tag	UNP Q46944
E	-3	GLY	-	expression tag	UNP Q46944
E	-2	SER	-	expression tag	UNP Q46944
E	-1	SER	-	expression tag	UNP Q46944
E	0	MET	-	expression tag	UNP Q46944
E	1	GLY	-	expression tag	UNP Q46944
E	69	LYS	LEU	engineered mutation	UNP Q46944
E	142	GLN	GLU	engineered mutation	UNP Q46944
F	-20	MET	-	expression tag	UNP Q46944
F	-19	ALA	-	expression tag	UNP Q46944
F	-18	SER	-	expression tag	UNP Q46944
F	-17	SER	-	expression tag	UNP Q46944
F	-16	HIS	-	expression tag	UNP Q46944
F	-15	HIS	-	expression tag	UNP Q46944
F	-14	HIS	-	expression tag	UNP Q46944
F	-13	HIS	-	expression tag	UNP Q46944
F	-12	HIS	-	expression tag	UNP Q46944
F	-11	HIS	-	expression tag	UNP Q46944
F	-10	SER	-	expression tag	UNP Q46944
F	-9	SER	-	expression tag	UNP Q46944
F	-8	GLY	-	expression tag	UNP Q46944
F	-7	LEU	-	expression tag	UNP Q46944
F	-6	VAL	-	expression tag	UNP Q46944
F	-5	PRO	-	expression tag	UNP Q46944
F	-4	ARG	-	expression tag	UNP Q46944
F	-3	GLY	-	expression tag	UNP Q46944
F	-2	SER	-	expression tag	UNP Q46944

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	SER	-	expression tag	UNP Q46944
F	0	MET	-	expression tag	UNP Q46944
F	1	GLY	-	expression tag	UNP Q46944
F	69	LYS	LEU	engineered mutation	UNP Q46944
F	142	GLN	GLU	engineered mutation	UNP Q46944
G	-20	MET	-	expression tag	UNP Q46944
G	-19	ALA	-	expression tag	UNP Q46944
G	-18	SER	-	expression tag	UNP Q46944
G	-17	SER	-	expression tag	UNP Q46944
G	-16	HIS	-	expression tag	UNP Q46944
G	-15	HIS	-	expression tag	UNP Q46944
G	-14	HIS	-	expression tag	UNP Q46944
G	-13	HIS	-	expression tag	UNP Q46944
G	-12	HIS	-	expression tag	UNP Q46944
G	-11	HIS	-	expression tag	UNP Q46944
G	-10	SER	-	expression tag	UNP Q46944
G	-9	SER	-	expression tag	UNP Q46944
G	-8	GLY	-	expression tag	UNP Q46944
G	-7	LEU	-	expression tag	UNP Q46944
G	-6	VAL	-	expression tag	UNP Q46944
G	-5	PRO	-	expression tag	UNP Q46944
G	-4	ARG	-	expression tag	UNP Q46944
G	-3	GLY	-	expression tag	UNP Q46944
G	-2	SER	-	expression tag	UNP Q46944
G	-1	SER	-	expression tag	UNP Q46944
G	0	MET	-	expression tag	UNP Q46944
G	1	GLY	-	expression tag	UNP Q46944
G	69	LYS	LEU	engineered mutation	UNP Q46944
G	142	GLN	GLU	engineered mutation	UNP Q46944
H	-20	MET	-	expression tag	UNP Q46944
H	-19	ALA	-	expression tag	UNP Q46944
H	-18	SER	-	expression tag	UNP Q46944
H	-17	SER	-	expression tag	UNP Q46944
H	-16	HIS	-	expression tag	UNP Q46944
H	-15	HIS	-	expression tag	UNP Q46944
H	-14	HIS	-	expression tag	UNP Q46944
H	-13	HIS	-	expression tag	UNP Q46944
H	-12	HIS	-	expression tag	UNP Q46944
H	-11	HIS	-	expression tag	UNP Q46944
H	-10	SER	-	expression tag	UNP Q46944
H	-9	SER	-	expression tag	UNP Q46944
H	-8	GLY	-	expression tag	UNP Q46944

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Chain	Residue	Modelled	Actual	Comment	Reference
H	-7	LEU	-	expression tag	UNP Q46944
H	-6	VAL	-	expression tag	UNP Q46944
H	-5	PRO	-	expression tag	UNP Q46944
H	-4	ARG	-	expression tag	UNP Q46944
H	-3	GLY	-	expression tag	UNP Q46944
H	-2	SER	-	expression tag	UNP Q46944
H	-1	SER	-	expression tag	UNP Q46944
H	0	MET	-	expression tag	UNP Q46944
H	1	GLY	-	expression tag	UNP Q46944
H	69	LYS	LEU	engineered mutation	UNP Q46944
H	142	GLN	GLU	engineered mutation	UNP Q46944

- Molecule 2 is a DNA chain called DNA (5'-D(P*CP*GP*GP*GP*AP*GP*GP*CP*CP*CP*GP*CP*GP*GP*GP*CP*CP*GP*CP*CP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	I	22	Total 453	C 210	N 90	O 131	P 22	0	0	0
2	K	22	Total 453	C 210	N 90	O 131	P 22	0	0	0
2	M	22	Total 453	C 210	N 90	O 131	P 22	0	0	0
2	O	22	Total 453	C 210	N 90	O 131	P 22	0	0	0

- Molecule 3 is a DNA chain called DNA (5'-D(P*GP*CP*GP*GP*CP*GP*GP*CP*CP*CP*GP*CP*GP*GP*GP*CP*CP*TP*CP*CP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	J	22	Total 449	C 209	N 85	O 133	P 22	0	0	0
3	L	22	Total 449	C 209	N 85	O 133	P 22	0	0	0
3	N	22	Total 449	C 209	N 85	O 133	P 22	0	0	0
3	P	22	Total 449	C 209	N 85	O 133	P 22	0	0	0

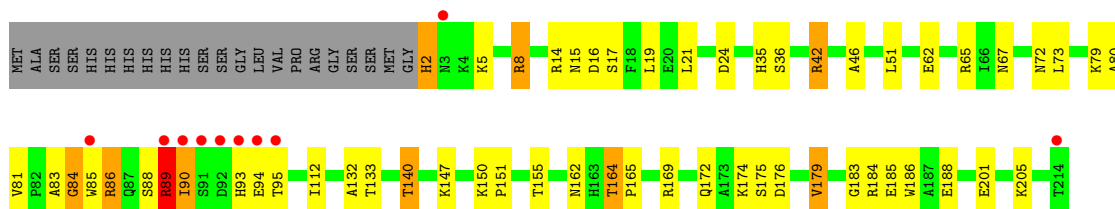
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	74	Total O 74 74	0	0
4	B	79	Total O 79 79	0	0
4	C	30	Total O 30 30	0	0
4	D	28	Total O 28 28	0	0
4	E	43	Total O 43 43	0	0
4	F	46	Total O 46 46	0	0
4	G	66	Total O 66 66	0	0
4	H	40	Total O 40 40	0	0
4	I	8	Total O 8 8	0	0
4	J	8	Total O 8 8	0	0
4	K	15	Total O 15 15	0	0
4	L	8	Total O 8 8	0	0
4	M	8	Total O 8 8	0	0
4	N	12	Total O 12 12	0	0
4	O	15	Total O 15 15	0	0
4	P	13	Total O 13 13	0	0

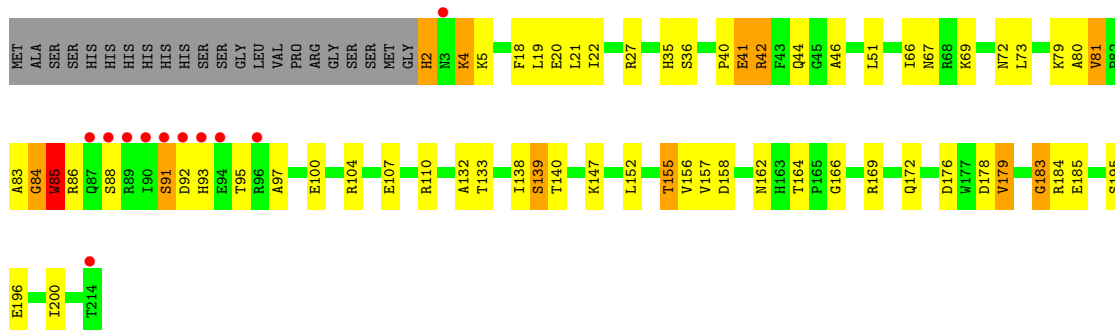
3 Residue-property plots [i](#)

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.

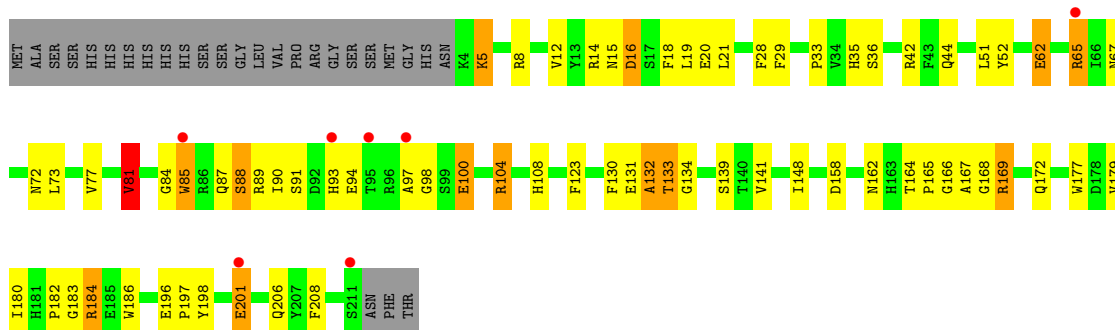
- Molecule 1: Eco29kIR



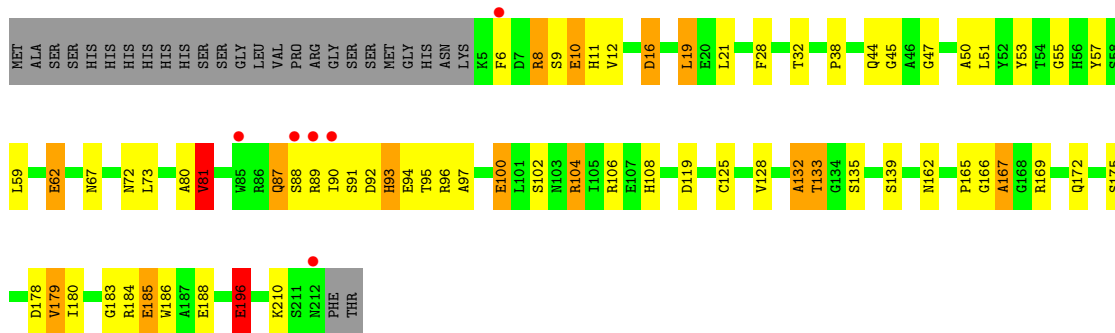
- Molecule 1: Eco29kIR



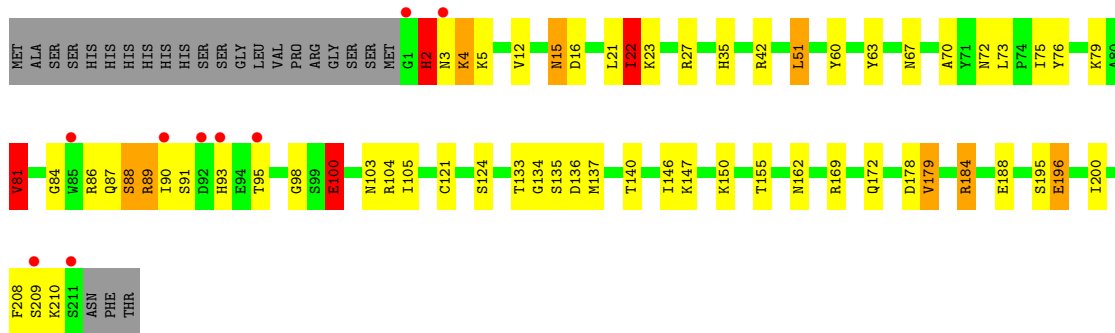
- Molecule 1: Eco29kIR



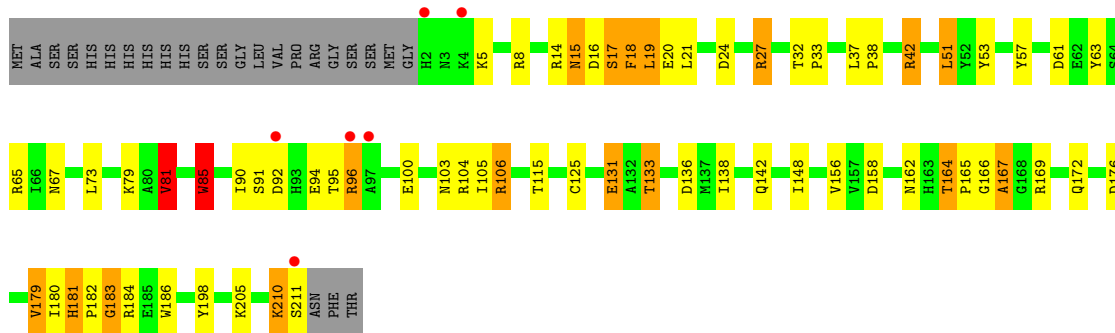
• Molecule 1: Eco29kIR



• Molecule 1: Eco29kIR

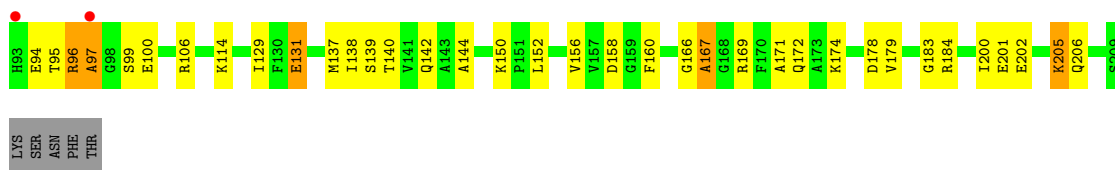


• Molecule 1: Eco29kIR

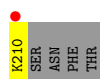
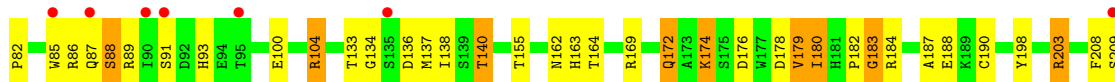


• Molecule 1: Eco29kIR





- Molecule 1: Eco29kIR



- Molecule 2: DNA (5'-D(P*CP*GP*GP*GP*AP*GP*GP*CP*CP*CP*GP*CP*GP*GP*GP*C P*CP*GP*CP*CP*GP*C)-3')



- Molecule 2: DNA (5'-D(P*CP*GP*GP*GP*AP*GP*GP*CP*CP*CP*GP*CP*GP*GP*GP*C P*CP*GP*CP*CP*GP*C)-3')



- Molecule 2: DNA (5'-D(P*CP*GP*GP*GP*AP*GP*GP*CP*CP*CP*GP*CP*GP*GP*GP*C P*CP*GP*CP*CP*GP*C)-3')



- Molecule 2: DNA (5'-D(P*CP*GP*GP*GP*AP*GP*GP*CP*CP*CP*GP*CP*GP*GP*GP*C P*CP*GP*CP*CP*GP*C)-3')





- Molecule 3: DNA (5'-D(P*GP*CP*GP*GP*CP*GP*GP*CP*CP*CP*GP*CP*GP*GP*GP*CP*CP*TP*CP*CP*CP*G)-3')



- Molecule 3: DNA (5'-D(P*GP*CP*GP*GP*CP*GP*GP*CP*CP*CP*GP*CP*GP*GP*GP*CP*CP*TP*CP*CP*CP*G)-3')



- Molecule 3: DNA (5'-D(P*GP*CP*GP*GP*CP*GP*GP*CP*CP*CP*GP*CP*GP*GP*GP*CP*CP*TP*CP*CP*CP*G)-3')



- Molecule 3: DNA (5'-D(P*GP*CP*GP*GP*CP*GP*GP*CP*CP*CP*GP*CP*GP*GP*GP*CP*CP*TP*CP*CP*CP*G)-3')



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	99.86Å 101.46Å 144.40Å 90.00° 110.19° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (50.00-2.50) 99.5 (50.00-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.28 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.211 , 0.274 0.219 , 0.275	Depositor DCC
R_{free} test set	4677 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	45.9	Xtrriage
Anisotropy	0.734	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17406	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 52.34 % 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 4.9318e-05. 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 i

5.1 Standard geometry i

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	1.21	0/1725	1.25	7/2334 (0.3%)
1	B	1.16	1/1738 (0.1%)	1.27	12/2357 (0.5%)
1	C	1.04	2/1698 (0.1%)	1.17	10/2302 (0.4%)
1	D	1.04	3/1682 (0.2%)	1.19	6/2279 (0.3%)
1	E	1.13	1/1713 (0.1%)	1.29	16/2325 (0.7%)
1	F	1.15	1/1723 (0.1%)	1.30	15/2333 (0.6%)
1	G	1.21	2/1712 (0.1%)	1.32	18/2322 (0.8%)
1	H	1.11	2/1708 (0.1%)	1.21	10/2317 (0.4%)
2	I	0.73	0/508	1.55	9/782 (1.2%)
2	K	0.69	0/508	1.49	7/782 (0.9%)
2	M	0.74	0/508	1.55	15/782 (1.9%)
2	O	0.73	0/508	1.60	13/782 (1.7%)
3	J	0.73	0/502	1.51	6/772 (0.8%)
3	L	0.67	0/502	1.70	14/772 (1.8%)
3	N	0.72	0/502	1.63	8/772 (1.0%)
3	P	0.73	0/502	1.65	14/772 (1.8%)
All	All	1.05	12/17739 (0.1%)	1.34	180/24785 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	F	0	1
1	G	0	2
1	H	0	2
All	All	0	7

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	87	GLN	CA-C	7.85	1.56	1.52
1	G	144	ALA	CA-CB	-7.67	1.41	1.53
1	H	180	ILE	CA-CB	6.56	1.62	1.54
1	G	90	ILE	CA-CB	5.97	1.62	1.54
1	E	12	VAL	CA-CB	5.76	1.60	1.54
1	D	45	GLY	C-O	5.62	1.28	1.23
1	C	133	THR	CA-CB	5.44	1.61	1.53
1	B	132	ALA	C-O	-5.44	1.19	1.24
1	H	174	LYS	CA-C	5.12	1.59	1.52
1	C	77	VAL	CA-CB	-5.05	1.48	1.54
1	D	104	ARG	CZ-NH2	-5.02	1.26	1.33
1	F	148	ILE	C-O	-5.01	1.18	1.24

All (180) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	85	TRP	N-CA-C	12.87	127.39	108.60
1	F	131	GLU	N-CA-C	11.09	123.12	111.14
1	G	131	GLU	N-CA-C	11.04	122.88	111.07
1	F	167	ALA	N-CA-C	-10.70	96.41	110.43
3	J	12	DC	O3'-P-O5'	-9.59	89.62	104.00
1	G	96	ARG	N-CA-C	9.29	122.79	110.53
1	H	81	VAL	CA-C-N	-8.96	110.80	119.85
1	H	81	VAL	C-N-CA	-8.96	110.80	119.85
1	E	150	LYS	CA-C-N	-8.48	111.28	119.85
1	E	150	LYS	C-N-CA	-8.48	111.28	119.85
3	P	16	DC	C4'-C3'-O3'	-8.36	97.46	110.00
1	B	81	VAL	CA-C-N	-8.16	111.60	119.85
1	B	81	VAL	C-N-CA	-8.16	111.60	119.85
3	N	8	DC	O3'-P-O5'	-8.06	91.92	104.00
3	L	10	DC	C2'-C3'-O3'	-7.92	99.62	111.50
1	A	94	GLU	N-CA-C	7.76	120.49	111.02
3	J	10	DC	C2'-C3'-O3'	-7.65	100.02	111.50
2	O	7	DG	C4'-C3'-O3'	-7.60	98.60	110.00
3	L	4	DG	P-O3'-C3'	7.26	131.09	120.20
2	K	15	DG	C4'-C3'-O3'	-7.08	99.38	110.00
1	H	47	GLY	N-CA-C	7.03	120.37	110.74
1	F	181	HIS	CA-C-N	-7.02	113.47	120.21
1	F	181	HIS	C-N-CA	-7.02	113.47	120.21
2	I	1	DC	C5'-C4'-C3'	6.96	125.34	114.90
1	F	81	VAL	CA-C-N	-6.93	112.64	119.78
1	F	81	VAL	C-N-CA	-6.93	112.64	119.78
1	E	22	ILE	N-CA-CB	6.87	120.88	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	N	20	DC	O3'-P-O5'	-6.86	93.70	104.00
3	P	4	DG	P-O3'-C3'	6.84	130.46	120.20
3	L	17	DC	C4'-C3'-O3'	-6.83	99.76	110.00
2	I	12	DC	C4'-C3'-O3'	-6.78	99.83	110.00
3	L	13	DG	O4'-C1'-N9	6.75	118.53	108.40
3	J	4	DG	P-O3'-C3'	6.68	130.22	120.20
1	H	28	PHE	N-CA-C	-6.68	103.61	111.03
2	K	14	DG	O4'-C1'-N9	-6.63	98.45	108.40
3	N	16	DC	C4'-C3'-O3'	-6.52	100.22	110.00
1	G	81	VAL	CA-C-N	-6.49	112.97	119.92
1	G	81	VAL	C-N-CA	-6.49	112.97	119.92
1	H	164	THR	CA-C-N	6.47	126.24	119.05
1	H	164	THR	C-N-CA	6.47	126.24	119.05
1	G	129	ILE	CB-CA-C	-6.47	103.13	110.96
1	F	92	ASP	N-CA-C	6.46	121.96	111.37
2	O	5	DA	O4'-C1'-N9	-6.39	98.82	108.40
2	O	7	DG	P-O3'-C3'	6.38	129.76	120.20
1	G	97	ALA	N-CA-C	6.36	119.87	111.28
3	N	12	DC	O3'-P-O5'	-6.33	94.50	104.00
1	G	150	LYS	CA-C-N	-6.26	113.33	119.78
1	G	150	LYS	C-N-CA	-6.26	113.33	119.78
1	E	196	GLU	CA-C-N	6.25	125.47	118.97
1	E	196	GLU	C-N-CA	6.25	125.47	118.97
1	B	41	GLU	N-CA-C	6.24	119.20	109.52
1	E	200	ILE	N-CA-C	-6.22	104.45	110.42
1	B	158	ASP	N-CA-C	6.21	118.80	110.35
1	G	4	LYS	N-CA-C	6.14	116.67	108.38
1	F	91	SER	N-CA-C	6.13	119.36	109.79
2	M	8	DC	P-O5'-C5'	6.13	129.20	120.00
2	O	6	DG	C2'-C3'-O3'	-6.11	102.33	111.50
1	G	158	ASP	CA-C-N	-6.07	116.28	123.08
1	G	158	ASP	C-N-CA	-6.07	116.28	123.08
1	A	112	ILE	N-CA-C	-6.06	104.40	111.00
1	E	179	VAL	N-CA-CB	6.02	118.28	110.57
1	E	22	ILE	CB-CA-C	-6.02	103.92	112.22
1	B	155	THR	N-CA-C	6.00	120.74	113.41
1	F	158	ASP	N-CA-C	5.97	118.62	110.55
1	E	195	SER	CA-C-N	5.97	128.67	120.67
1	E	195	SER	C-N-CA	5.97	128.67	120.67
2	K	17	DC	C4'-C3'-O3'	-5.97	101.05	110.00
1	E	81	VAL	CA-C-N	-5.95	113.81	119.76
1	E	81	VAL	C-N-CA	-5.95	113.81	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	12	DC	N1-C1'-C2'	5.95	122.42	113.50
3	L	5	DC	O3'-P-O5'	-5.93	95.10	104.00
2	O	3	DG	O5'-C5'-C4'	-5.92	101.92	110.80
1	E	75	ILE	N-CA-C	-5.92	104.69	113.39
2	I	13	DG	O4'-C1'-N9	5.92	117.28	108.40
3	P	20	DC	C4'-C3'-O3'	-5.92	101.12	110.00
3	N	15	DG	P-O5'-C5'	5.92	128.87	120.00
1	H	91	SER	N-CA-C	5.91	118.82	110.14
3	L	6	DG	O4'-C1'-N9	-5.88	99.58	108.40
3	L	6	DG	C4'-C3'-O3'	-5.87	101.20	110.00
1	C	85	TRP	N-CA-C	5.84	118.43	111.71
1	G	167	ALA	N-CA-C	-5.82	98.40	110.80
2	M	13	DG	C5'-C4'-C3'	-5.82	106.16	114.90
1	A	184	ARG	N-CA-C	-5.82	101.24	109.96
1	C	148	ILE	N-CA-C	5.80	115.98	110.53
2	M	6	DG	C4'-C3'-O3'	-5.78	101.33	110.00
2	M	12	DC	O3'-P-O5'	-5.78	95.33	104.00
1	G	138	ILE	N-CA-C	5.77	116.42	110.36
2	M	5	DA	C4'-C3'-O3'	-5.76	101.36	110.00
1	A	155	THR	CA-C-N	-5.76	117.06	123.16
1	A	155	THR	C-N-CA	-5.76	117.06	123.16
2	M	8	DC	N1-C1'-C2'	5.74	122.11	113.50
3	N	10	DC	P-O3'-C3'	5.73	128.80	120.20
2	K	13	DG	O4'-C1'-N9	5.71	116.97	108.40
1	C	196	GLU	CA-C-N	5.68	126.94	119.84
1	C	196	GLU	C-N-CA	5.68	126.94	119.84
1	F	105	ILE	N-CA-CB	5.65	117.16	110.55
1	C	98	GLY	N-CA-C	5.65	117.97	111.36
1	D	167	ALA	N-CA-C	-5.64	98.79	110.80
3	P	13	DG	O3'-P-O5'	-5.63	95.55	104.00
2	O	4	DG	C2'-C3'-O3'	-5.63	103.05	111.50
3	L	7	DG	C4'-C3'-O3'	-5.63	101.56	110.00
3	L	17	DC	P-O3'-C3'	5.62	128.63	120.20
3	L	7	DG	P-O3'-C3'	5.62	128.62	120.20
3	P	15	DG	C4'-C3'-O3'	-5.60	101.60	110.00
1	D	38	PRO	CA-C-N	5.59	123.74	119.66
1	D	38	PRO	C-N-CA	5.59	123.74	119.66
3	P	11	DG	O4'-C1'-N9	-5.58	100.03	108.40
2	I	13	DG	C2'-C3'-O3'	-5.58	103.14	111.50
3	J	11	DG	O4'-C1'-N9	-5.56	100.06	108.40
1	F	156	VAL	CB-CA-C	-5.53	106.13	111.05
1	A	132	ALA	CB-CA-C	-5.53	110.22	116.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	184	ARG	N-CA-C	-5.52	100.98	109.76
2	M	16	DC	O3'-P-O5'	-5.52	95.73	104.00
1	G	85	TRP	N-CA-C	5.51	117.75	111.02
1	C	81	VAL	N-CA-CB	-5.51	103.49	111.21
2	M	10	DC	P-O3'-C3'	5.51	128.46	120.20
1	E	100	GLU	N-CA-C	5.50	117.27	111.28
1	H	88	SER	N-CA-C	5.46	122.42	110.80
2	K	14	DG	N9-C1'-C2'	5.43	121.64	113.50
3	J	4	DG	C2'-C3'-O3'	5.42	119.64	111.50
3	P	6	DG	O5'-C5'-C4'	5.42	118.94	110.80
1	C	20	GLU	N-CA-C	-5.42	105.46	111.36
2	I	6	DG	O4'-C1'-N9	-5.40	100.30	108.40
3	P	20	DC	C2'-C3'-O3'	5.40	119.59	111.50
3	P	9	DC	N1-C1'-C2'	5.39	121.59	113.50
2	M	13	DG	O4'-C1'-N9	5.38	116.47	108.40
2	O	13	DG	C2'-C3'-O3'	-5.36	103.45	111.50
1	F	164	THR	CA-C-N	5.36	125.17	119.28
1	F	164	THR	C-N-CA	5.36	125.17	119.28
3	J	10	DC	O3'-P-O5'	-5.36	95.96	104.00
2	M	6	DG	P-O3'-C3'	5.32	128.18	120.20
2	I	1	DC	O5'-C5'-C4'	5.32	118.78	110.80
1	H	32	THR	CA-C-N	5.30	125.61	119.93
1	H	32	THR	C-N-CA	5.30	125.61	119.93
2	O	17	DC	C4'-C3'-O3'	-5.30	102.05	110.00
1	C	169	ARG	N-CA-C	5.29	119.28	112.41
1	B	69	LYS	N-CA-C	-5.29	107.00	113.50
2	I	1	DC	P-O3'-C3'	5.28	128.12	120.20
3	P	18	DT	C4'-C3'-O3'	-5.27	102.09	110.00
1	D	196	GLU	CA-C-N	5.27	126.43	119.84
1	D	196	GLU	C-N-CA	5.27	126.43	119.84
2	O	8	DC	O3'-P-O5'	-5.26	96.11	104.00
3	L	18	DT	C4'-C3'-O3'	-5.24	102.15	110.00
3	L	13	DG	P-O3'-C3'	5.23	128.05	120.20
3	N	13	DG	O4'-C1'-N9	5.23	116.24	108.40
3	L	17	DC	O3'-P-O5'	-5.22	96.18	104.00
1	D	81	VAL	N-CA-CB	-5.21	103.91	111.21
3	P	4	DG	C2'-C3'-O3'	5.21	119.31	111.50
1	G	200	ILE	CB-CA-C	-5.21	105.20	112.02
1	G	142	GLN	N-CA-C	-5.20	105.61	111.28
1	E	184	ARG	CB-CG-CD	-5.19	99.36	111.30
1	C	134	GLY	N-CA-C	-5.19	100.88	113.18
2	I	3	DG	P-O3'-C3'	5.19	127.98	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	105	ILE	N-CA-C	-5.19	105.33	110.62
2	M	15	DG	N9-C1'-C2'	5.19	121.28	113.50
3	N	4	DG	P-O3'-C3'	5.18	127.97	120.20
2	O	16	DC	N1-C1'-C2'	5.17	121.25	113.50
2	I	10	DC	P-O3'-C3'	5.16	127.94	120.20
2	M	16	DC	C4'-C3'-O3'	-5.16	102.27	110.00
1	B	88	SER	N-CA-C	5.16	119.04	112.34
3	P	13	DG	P-O3'-C3'	5.16	127.93	120.20
2	K	17	DC	P-O3'-C3'	5.14	127.91	120.20
1	B	4	LYS	N-CA-C	5.12	121.70	110.80
1	F	85	TRP	N-CA-C	5.12	117.25	111.11
2	K	14	DG	O3'-P-O5'	5.11	111.67	104.00
1	B	138	ILE	N-CA-C	5.11	115.26	110.30
2	M	10	DC	C2'-C3'-O3'	-5.10	103.85	111.50
1	F	138	ILE	N-CA-C	5.08	115.51	110.23
2	M	7	DG	N9-C1'-C2'	5.07	121.10	113.50
2	M	13	DG	O3'-P-O5'	-5.06	96.41	104.00
2	O	10	DC	C2'-C3'-O3'	-5.06	103.91	111.50
1	A	5	LYS	N-CA-C	5.05	116.81	110.24
1	B	81	VAL	N-CA-CB	-5.05	104.14	111.21
1	B	139	SER	N-CA-C	5.05	117.44	111.33
1	G	26	VAL	CA-C-N	5.04	126.99	120.44
1	G	26	VAL	C-N-CA	5.04	126.99	120.44
3	L	11	DG	P-O5'-C5'	5.04	127.55	120.00
2	O	7	DG	O5'-C5'-C4'	-5.03	103.25	110.80
3	P	20	DC	N1-C1'-C2'	5.03	121.04	113.50
3	P	11	DG	N9-C1'-C2'	5.01	121.02	113.50

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	85	TRP	Peptide
1	B	183	GLY	Peptide
1	F	183	GLY	Peptide
1	G	183	GLY	Peptide
1	G	96	ARG	Peptide
1	H	183	GLY	Peptide
1	H	5	LYS	Peptide

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	1677	0	1578	51	1
1	B	1687	0	1567	68	0
1	C	1649	0	1540	59	0
1	D	1634	0	1524	57	0
1	E	1663	0	1544	59	0
1	F	1674	0	1578	58	0
1	G	1663	0	1574	41	1
1	H	1658	0	1544	55	0
2	I	453	0	243	4	0
2	K	453	0	243	8	0
2	M	453	0	243	12	1
2	O	453	0	243	15	1
3	J	449	0	244	7	0
3	L	449	0	244	11	0
3	N	449	0	244	14	0
3	P	449	0	244	11	0
4	A	74	0	0	4	0
4	B	79	0	0	7	0
4	C	30	0	0	1	0
4	D	28	0	0	3	0
4	E	43	0	0	8	0
4	F	46	0	0	4	0
4	G	66	0	0	0	0
4	H	40	0	0	1	0
4	I	8	0	0	0	0
4	J	8	0	0	0	0
4	K	15	0	0	1	0
4	L	8	0	0	0	0
4	M	8	0	0	0	0
4	N	12	0	0	0	0
4	O	15	0	0	2	0
4	P	13	0	0	4	0
All	All	17406	0	14397	411	2

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 13.

All (411) 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:E:2:HIS:HB2	1:E:3:ASN:HA	1.18	1.17
1:E:121:CYS:HB2	4:E:408:HOH:O	1.45	1.15
1:C:87:GLN:HB3	1:C:88:SER:CA	1.82	1.09
1:A:140:THR:HG21	1:H:140:THR:HG22	1.39	1.05
1:H:104:ARG:HH11	1:H:104:ARG:HG2	1.21	1.03
1:B:85:TRP:HE3	1:B:85:TRP:HA	1.21	1.03
1:B:85:TRP:HA	1:B:85:TRP:CE3	1.90	1.02
1:D:87:GLN:HG3	1:D:88:SER:H	1.20	1.01
1:G:81:VAL:HG11	1:G:85:TRP:CD1	1.97	1.00
1:A:140:THR:HG21	1:H:140:THR:CG2	1.92	1.00
1:B:86:ARG:HG2	4:P:415:HOH:O	1.62	0.97
3:N:1:DG:H1'	3:N:2:DC:C5	2.02	0.94
1:F:106:ARG:HG2	1:F:106:ARG:HH11	1.31	0.93
1:B:172:GLN:HE22	1:E:169:ARG:HH21	1.10	0.93
1:B:67:ASN:HD21	1:B:73:LEU:H	1.08	0.93
1:F:67:ASN:HD21	1:F:73:LEU:H	1.11	0.93
1:B:107:GLU:HG2	4:B:321:HOH:O	1.68	0.92
1:G:68:ARG:HE	1:G:68:ARG:HA	1.34	0.91
1:G:67:ASN:HD21	1:G:73:LEU:H	1.20	0.89
1:H:27:ARG:HG3	1:H:27:ARG:HH11	1.38	0.89
1:D:104:ARG:O	1:D:108:HIS:HD2	1.55	0.88
1:H:203:ARG:HG3	1:H:203:ARG:HH11	1.39	0.87
1:F:104:ARG:HG3	1:F:104:ARG:HH11	1.40	0.86
1:A:147:LYS:HE3	4:A:241:HOH:O	1.75	0.85
1:E:87:GLN:HB3	1:E:88:SER:CA	2.07	0.84
1:A:140:THR:HB	1:H:136:ASP:O	1.77	0.84
1:D:87:GLN:CG	1:D:88:SER:H	1.91	0.83
1:E:2:HIS:HB2	1:E:3:ASN:CA	2.05	0.82
1:H:104:ARG:HH11	1:H:104:ARG:CG	1.90	0.82
1:F:180:ILE:HD12	1:F:181:HIS:CD2	2.14	0.82
1:C:33:PRO:HB3	1:G:4:LYS:HE2	1.63	0.81
1:B:104:ARG:HD3	4:B:428:HOH:O	1.81	0.79
1:A:67:ASN:ND2	1:A:72:ASN:H	1.81	0.79
2:O:1:DC:O2	3:P:22:DG:N2	2.16	0.79
1:E:22:ILE:HD12	1:E:23:LYS:H	1.46	0.78
3:P:1:DG:H4'	3:P:2:DC:OP1	1.82	0.78
1:C:172:GLN:NE2	1:G:169:ARG:HE	1.81	0.77
1:C:172:GLN:HE22	1:G:169:ARG:HH21	1.32	0.77
1:B:169:ARG:HH21	1:E:172:GLN:HE22	1.30	0.76
1:E:67:ASN:HD21	1:E:73:LEU:H	1.32	0.76
3:L:21:DC:H2''	3:L:22:DG:O5'	1.85	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:GLN:HE22	1:H:169:ARG:HH21	1.32	0.76
3:P:21:DC:H5	4:P:311:HOH:O	1.68	0.76
1:B:162:ASN:HD21	3:P:11:DG:H2'	1.49	0.76
1:B:42:ARG:HH12	2:O:5:DA:H5''	1.51	0.75
1:B:83:ALA:O	1:B:84:GLY:C	2.29	0.75
1:C:172:GLN:HE21	1:G:169:ARG:HE	1.35	0.74
1:E:2:HIS:CB	1:E:3:ASN:HA	2.09	0.74
2:K:1:DC:O2	3:L:22:DG:N2	2.18	0.73
1:B:172:GLN:HE22	1:E:169:ARG:NH2	1.84	0.73
1:A:67:ASN:HD21	1:A:73:LEU:H	1.38	0.72
1:F:42:ARG:HD3	1:F:106:ARG:HE	1.55	0.72
1:B:42:ARG:HH12	2:O:5:DA:C5'	2.02	0.72
3:P:3:DG:H2''	3:P:4:DG:C8	2.25	0.71
1:B:85:TRP:CE3	1:B:85:TRP:CA	2.73	0.71
1:D:96:ARG:N	1:D:97:ALA:CA	2.54	0.71
1:C:65:ARG:HG2	1:C:65:ARG:HH11	1.57	0.70
1:B:2:HIS:CD2	1:E:35:HIS:HD2	2.10	0.69
1:H:67:ASN:HD21	1:H:73:LEU:H	1.39	0.69
1:C:87:GLN:HB3	1:C:88:SER:C	2.16	0.69
1:A:2:HIS:HB3	1:H:35:HIS:CD2	2.27	0.69
1:E:22:ILE:HD12	1:E:23:LYS:N	2.07	0.69
2:O:5:DA:N3	4:O:350:HOH:O	2.25	0.69
1:B:172:GLN:NE2	1:E:169:ARG:HE	1.91	0.68
1:C:67:ASN:HD21	1:C:73:LEU:H	1.40	0.68
1:F:95:THR:HA	1:F:96:ARG:CB	2.23	0.68
1:H:104:ARG:CG	1:H:104:ARG:NH1	2.51	0.68
2:M:22:DC:O2	3:N:1:DG:N2	2.20	0.68
1:H:203:ARG:HH11	1:H:203:ARG:CG	2.08	0.67
1:G:81:VAL:CG1	1:G:85:TRP:CD1	2.75	0.67
1:A:67:ASN:HD22	1:A:72:ASN:H	1.43	0.67
1:D:67:ASN:HD21	1:D:73:LEU:H	1.41	0.66
2:M:18:DG:H1	3:N:5:DC:H42	1.41	0.66
1:C:172:GLN:NE2	1:G:169:ARG:HH21	1.93	0.66
1:D:169:ARG:HH21	1:F:172:GLN:HE22	1.44	0.66
1:B:85:TRP:CD1	1:B:86:ARG:HE	2.14	0.66
1:D:172:GLN:NE2	1:F:169:ARG:HE	1.93	0.66
1:B:2:HIS:HD2	1:E:35:HIS:HD2	1.42	0.65
1:E:67:ASN:ND2	1:E:72:ASN:H	1.93	0.65
1:B:183:GLY:HA2	4:B:352:HOH:O	1.96	0.65
3:P:21:DC:H2''	3:P:22:DG:O5'	1.95	0.65
1:D:12:VAL:HG13	1:F:131:GLU:HG3	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:64:SER:O	1:H:68:ARG:NH1	2.30	0.65
1:D:19:LEU:HD23	1:F:133:THR:HG22	1.79	0.64
1:C:139:SER:HB2	1:G:85:TRP:HH2	1.62	0.64
1:D:87:GLN:HG3	1:D:88:SER:N	2.03	0.64
1:E:84:GLY:O	1:E:88:SER:CA	2.45	0.64
1:D:166:GLY:C	1:D:167:ALA:O	2.40	0.64
1:H:27:ARG:HG3	1:H:27:ARG:NH1	2.10	0.64
1:C:172:GLN:HE22	1:G:169:ARG:NH2	1.95	0.63
1:A:183:GLY:HA2	4:A:314:HOH:O	1.97	0.63
1:D:87:GLN:CG	1:D:88:SER:N	2.61	0.63
1:D:169:ARG:HE	1:F:172:GLN:NE2	1.95	0.63
3:L:3:DG:H2''	3:L:4:DG:C8	2.33	0.63
3:N:2:DC:H2''	3:N:3:DG:C8	2.34	0.63
2:I:1:DC:O2	3:J:22:DG:N2	2.17	0.63
1:H:104:ARG:HG2	1:H:104:ARG:NH1	2.01	0.63
1:C:84:GLY:HA3	1:C:89:ARG:CB	2.29	0.63
1:F:15:ASN:ND2	1:F:17:SER:HB3	2.12	0.63
1:D:183:GLY:HA2	4:D:355:HOH:O	1.98	0.62
1:F:180:ILE:CD1	1:F:181:HIS:CD2	2.82	0.62
1:D:169:ARG:HE	1:F:172:GLN:HE21	1.45	0.62
1:B:147:LYS:HE3	4:E:490:HOH:O	1.98	0.62
1:F:104:ARG:HG3	1:F:104:ARG:NH1	2.14	0.62
1:B:164:THR:HG22	1:B:166:GLY:H	1.64	0.62
1:A:162:ASN:HD22	2:K:12:DC:H5	1.47	0.62
1:B:162:ASN:ND2	3:P:11:DG:H2'	2.14	0.62
1:D:104:ARG:O	1:D:108:HIS:CD2	2.47	0.62
1:A:79:LYS:HE3	4:A:348:HOH:O	1.98	0.61
1:F:115:THR:HA	1:F:182:PRO:O	2.00	0.61
1:A:140:THR:CG2	1:H:140:THR:HG22	2.23	0.61
1:D:44:GLN:O	1:F:8:ARG:NH1	2.33	0.61
1:F:104:ARG:HH11	1:F:104:ARG:CG	2.13	0.61
1:B:91:SER:CB	1:F:57:TYR:HE1	2.13	0.61
1:F:106:ARG:HH11	1:F:106:ARG:CG	2.10	0.60
1:A:140:THR:CG2	1:H:140:THR:CG2	2.76	0.60
1:B:85:TRP:HZ2	1:E:146:ILE:HG22	1.65	0.60
1:D:185:GLU:O	1:D:188:GLU:HB2	2.01	0.60
1:C:89:ARG:O	1:C:91:SER:N	2.34	0.59
1:D:8:ARG:HD3	1:D:9:SER:H	1.66	0.59
2:K:18:DG:H1'	2:K:19:DC:H5'	1.83	0.59
1:A:2:HIS:HA	1:H:34:VAL:O	2.02	0.59
1:D:6:PHE:O	1:D:6:PHE:CD2	2.55	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:GLY:O	1:B:85:TRP:HB2	2.03	0.59
1:D:172:GLN:HE22	1:F:169:ARG:HH21	1.49	0.59
4:B:297:HOH:O	1:E:35:HIS:HE1	1.86	0.58
1:C:94:GLU:OE2	1:C:97:ALA:HB2	2.02	0.58
1:B:42:ARG:NH1	2:O:5:DA:H5''	2.16	0.58
1:E:86:ARG:C	4:E:327:HOH:O	2.45	0.58
1:A:83:ALA:O	1:A:84:GLY:C	2.47	0.58
1:C:16:ASP:OD1	1:C:16:ASP:N	2.37	0.58
1:B:85:TRP:HB3	4:P:415:HOH:O	2.02	0.58
1:D:80:ALA:HB2	1:D:104:ARG:HG3	1.85	0.58
1:G:44:GLN:OE1	1:G:97:ALA:HB1	2.03	0.58
1:B:40:PRO:HD3	1:E:2:HIS:CE1	2.39	0.58
1:D:132:ALA:O	1:D:133:THR:OG1	2.11	0.57
1:A:35:HIS:HE1	4:H:332:HOH:O	1.87	0.57
1:A:164:THR:HG23	1:A:186:TRP:CG	2.39	0.57
1:A:140:THR:CB	1:H:136:ASP:O	2.50	0.57
1:D:8:ARG:CD	1:D:9:SER:H	2.18	0.57
1:G:178:ASP:OD1	1:G:184:ARG:HD2	2.05	0.57
1:C:169:ARG:O	1:C:172:GLN:HG2	2.05	0.57
1:B:18:PHE:HZ	1:E:140:THR:HG21	1.70	0.57
1:D:8:ARG:HB3	1:D:10:GLU:HG2	1.87	0.57
1:D:19:LEU:HD13	1:F:136:ASP:OD2	2.05	0.57
1:B:169:ARG:O	1:B:172:GLN:HG2	2.05	0.57
1:C:81:VAL:HG13	1:C:85:TRP:HB2	1.87	0.57
1:F:24:ASP:HA	1:F:27:ARG:HH11	1.70	0.57
1:A:175:SER:O	1:A:179:VAL:HG12	2.05	0.56
1:E:81:VAL:O	1:E:100:GLU:HG3	2.04	0.56
1:B:35:HIS:HE1	4:E:286:HOH:O	1.88	0.56
1:F:211:SER:C	4:F:283:HOH:O	2.47	0.56
1:C:12:VAL:HG13	1:G:131:GLU:HG2	1.87	0.56
1:C:169:ARG:HH21	1:G:172:GLN:HE22	1.52	0.56
1:B:86:ARG:HD2	1:E:155:THR:HA	1.87	0.56
1:B:67:ASN:ND2	1:B:72:ASN:H	2.05	0.55
1:E:86:ARG:HA	4:E:472:HOH:O	2.06	0.55
1:E:162:ASN:HD22	2:O:12:DC:H5	1.53	0.55
1:B:2:HIS:HD2	1:E:35:HIS:CD2	2.22	0.55
1:F:16:ASP:O	1:F:19:LEU:HD22	2.06	0.55
3:P:2:DC:H2''	3:P:3:DG:C8	2.41	0.55
1:C:100:GLU:CD	1:C:100:GLU:H	2.14	0.55
1:D:172:GLN:HE21	1:F:169:ARG:HE	1.53	0.55
3:L:2:DC:H2''	3:L:3:DG:C8	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:162:ASN:ND2	3:L:12:DC:H5	2.05	0.55
1:B:19:LEU:HD21	1:E:137:MET:HE2	1.89	0.54
1:C:132:ALA:O	1:C:133:THR:OG1	2.16	0.54
1:A:21:LEU:C	1:A:21:LEU:HD13	2.32	0.54
1:A:176:ASP:HA	1:A:179:VAL:HG13	1.89	0.54
2:O:20:DC:H2''	2:O:21:DG:C8	2.43	0.54
1:C:198:TYR:O	1:C:201:GLU:HG3	2.07	0.54
1:B:85:TRP:CG	1:B:86:ARG:H	2.25	0.54
1:B:169:ARG:HH21	1:E:172:GLN:NE2	2.05	0.54
1:C:104:ARG:O	1:C:108:HIS:HD2	1.90	0.54
1:C:36:SER:N	1:G:2:HIS:HA	2.23	0.54
1:B:157:VAL:O	1:B:157:VAL:HG12	2.09	0.53
1:B:85:TRP:CH2	1:E:147:LYS:HA	2.44	0.53
1:H:81:VAL:HG13	1:H:82:PRO:HD2	1.91	0.53
1:C:65:ARG:HH11	1:C:65:ARG:CG	2.21	0.53
1:E:79:LYS:HE2	4:E:420:HOH:O	2.08	0.53
1:F:103:ASN:ND2	4:F:339:HOH:O	2.41	0.53
1:H:86:ARG:O	1:H:86:ARG:HG3	2.08	0.53
2:M:4:DG:H1	3:N:19:DC:H42	1.56	0.53
1:A:172:GLN:NE2	1:H:169:ARG:HE	2.06	0.53
1:F:104:ARG:HD2	4:F:432:HOH:O	2.09	0.53
1:G:160:PHE:O	1:G:184:ARG:NH2	2.41	0.53
1:C:67:ASN:ND2	1:C:72:ASN:H	2.06	0.53
2:M:22:DC:N3	3:N:1:DG:N1	2.38	0.53
1:C:162:ASN:HD22	2:M:12:DC:H5	1.56	0.53
1:D:44:GLN:HG2	1:D:97:ALA:O	2.09	0.53
2:K:12:DC:H2''	2:K:13:DG:O5'	2.08	0.53
1:B:169:ARG:HE	1:E:172:GLN:NE2	2.06	0.52
1:F:51:LEU:HD12	1:F:125:CYS:HB3	1.91	0.52
1:B:178:ASP:OD1	1:B:184:ARG:HD2	2.09	0.52
1:C:164:THR:HG22	1:C:166:GLY:H	1.75	0.52
1:G:152:LEU:HD12	1:G:156:VAL:HB	1.91	0.52
1:D:57:TYR:CE2	1:D:59:LEU:HB2	2.44	0.52
2:O:2:DG:H2''	2:O:3:DG:OP2	2.08	0.52
1:B:85:TRP:N	4:B:334:HOH:O	2.42	0.52
1:C:42:ARG:NH1	3:N:5:DC:OP1	2.43	0.52
1:E:16:ASP:HB3	1:F:65:ARG:HG3	1.92	0.52
1:A:169:ARG:HH21	1:H:172:GLN:HE22	1.58	0.52
1:F:32:THR:HB	1:F:33:PRO:HD2	1.92	0.52
1:B:84:GLY:O	1:B:85:TRP:CB	2.58	0.52
1:C:28:PHE:O	1:C:29:PHE:C	2.51	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:15:ASN:HD21	1:F:17:SER:HB3	1.73	0.51
1:D:162:ASN:HD22	3:J:12:DC:H5	1.59	0.51
1:A:19:LEU:HD21	1:H:137:MET:CE	2.41	0.51
1:B:42:ARG:NH1	2:O:5:DA:C5'	2.73	0.51
2:M:20:DC:H2''	2:M:21:DG:OP2	2.11	0.51
1:F:166:GLY:C	1:F:167:ALA:O	2.48	0.51
1:D:178:ASP:OD1	1:D:184:ARG:HD2	2.10	0.51
1:G:166:GLY:C	1:G:167:ALA:O	2.49	0.51
1:H:203:ARG:HG3	1:H:203:ARG:NH1	2.19	0.51
1:C:158:ASP:HA	4:C:379:HOH:O	2.11	0.51
1:E:178:ASP:OD1	1:E:184:ARG:HD2	2.11	0.51
2:K:20:DC:H3'	4:K:380:HOH:O	2.11	0.51
1:E:196:GLU:OE2	1:E:196:GLU:HA	2.11	0.50
1:H:176:ASP:HA	1:H:179:VAL:HG13	1.93	0.50
1:A:67:ASN:ND2	1:A:72:ASN:N	2.56	0.50
1:D:90:ILE:N	1:D:91:SER:CA	2.74	0.50
1:A:88:SER:O	1:A:90:ILE:N	2.45	0.50
2:O:6:DG:N3	4:O:342:HOH:O	2.34	0.50
1:E:86:ARG:HA	1:E:86:ARG:HE	1.76	0.50
1:E:15:ASN:C	1:E:15:ASN:HD22	2.18	0.50
2:I:16:DC:H1'	2:I:17:DC:H5'	1.94	0.50
1:B:67:ASN:HD22	1:B:72:ASN:H	1.60	0.50
1:B:79:LYS:HE3	4:P:329:HOH:O	2.11	0.50
1:D:102:SER:O	1:D:106:ARG:HG3	2.11	0.50
1:H:27:ARG:HH11	1:H:27:ARG:CG	2.17	0.49
1:C:18:PHE:CG	1:G:137:MET:HG3	2.47	0.49
1:C:169:ARG:HE	1:G:172:GLN:HE21	1.60	0.49
1:D:53:TYR:CZ	1:D:55:GLY:HA3	2.47	0.49
1:A:172:GLN:HE21	1:H:169:ARG:HE	1.59	0.49
2:M:18:DG:H1	3:N:5:DC:N4	2.10	0.49
1:B:183:GLY:CA	4:B:352:HOH:O	2.58	0.49
1:D:175:SER:O	1:D:179:VAL:HG12	2.13	0.49
1:A:183:GLY:HA3	4:A:394:HOH:O	2.12	0.49
1:C:172:GLN:NE2	1:G:169:ARG:NE	2.57	0.49
1:A:42:ARG:HH22	3:L:5:DC:H5''	1.78	0.49
1:F:42:ARG:HD3	1:F:106:ARG:NE	2.25	0.49
1:D:9:SER:C	1:D:11:HIS:H	2.21	0.49
1:D:172:GLN:NE2	1:F:169:ARG:HH21	2.10	0.49
1:H:67:ASN:ND2	1:H:72:ASN:H	2.11	0.49
1:B:172:GLN:HE21	1:E:169:ARG:HE	1.61	0.49
1:C:169:ARG:HH21	1:G:172:GLN:NE2	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:87:GLN:CB	1:C:88:SER:CA	2.69	0.48
1:D:81:VAL:O	1:D:100:GLU:HG3	2.11	0.48
1:F:106:ARG:HG2	1:F:106:ARG:NH1	2.08	0.48
1:G:16:ASP:O	1:G:19:LEU:HD12	2.12	0.48
1:C:197:PRO:O	1:C:198:TYR:C	2.53	0.48
1:E:86:ARG:HG2	2:O:15:DG:OP1	2.13	0.48
1:F:95:THR:CA	1:F:96:ARG:CB	2.90	0.48
2:O:1:DC:N3	3:P:22:DG:N1	2.49	0.48
1:D:92:ASP:CA	1:D:93:HIS:C	2.87	0.48
1:C:42:ARG:NH2	3:N:5:DC:OP1	2.46	0.48
1:D:94:GLU:N	1:D:95:THR:CA	2.77	0.48
1:C:15:ASN:HB3	1:G:137:MET:HE1	1.95	0.48
1:C:177:TRP:HZ3	1:C:184:ARG:HD2	1.78	0.48
3:J:1:DG:H1'	3:J:2:DC:C5	2.49	0.48
1:C:81:VAL:O	1:C:100:GLU:HG3	2.14	0.48
1:G:169:ARG:NH2	2:M:10:DC:H2''	2.29	0.48
1:F:81:VAL:HG11	1:F:85:TRP:CD1	2.49	0.47
1:A:36:SER:HB3	1:H:2:HIS:CD2	2.49	0.47
1:F:16:ASP:O	1:F:18:PHE:N	2.47	0.47
1:C:166:GLY:C	1:C:167:ALA:O	2.57	0.47
1:B:2:HIS:HA	1:E:35:HIS:HA	1.96	0.47
1:A:89:ARG:HE	1:A:89:ARG:HA	1.78	0.47
3:N:1:DG:H4'	3:N:2:DC:OP1	2.14	0.47
1:B:19:LEU:HD21	1:E:134:GLY:HA2	1.95	0.47
1:C:139:SER:HB2	1:G:85:TRP:CH2	2.46	0.47
1:F:104:ARG:NH1	1:F:104:ARG:CG	2.73	0.47
1:A:15:ASN:OD1	1:A:17:SER:HB2	2.14	0.47
1:B:83:ALA:O	1:B:84:GLY:O	2.31	0.47
1:H:134:GLY:O	1:H:137:MET:HB2	2.15	0.47
1:A:162:ASN:HD21	2:K:11:DG:H2'	1.79	0.47
1:C:130:PHE:CZ	1:C:141:VAL:HG21	2.50	0.47
1:F:183:GLY:HA2	4:F:306:HOH:O	2.15	0.47
1:G:19:LEU:HD12	1:G:19:LEU:H	1.79	0.47
1:H:27:ARG:HD3	1:H:27:ARG:N	2.30	0.47
1:B:133:THR:OG1	1:F:61:ASP:OD1	2.20	0.47
1:B:44:GLN:NE2	1:B:97:ALA:HB3	2.29	0.46
1:C:62:GLU:HG3	1:C:208:PHE:CG	2.50	0.46
1:B:139:SER:HB2	4:B:288:HOH:O	2.16	0.46
1:E:103:ASN:ND2	4:E:365:HOH:O	2.48	0.46
2:M:1:DC:O2	3:N:22:DG:N2	2.32	0.46
2:K:1:DC:N3	3:L:22:DG:N1	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:133:THR:O	1:G:18:PHE:HE1	1.98	0.46
2:I:17:DC:H2''	2:I:18:DG:C8	2.50	0.46
1:D:67:ASN:ND2	1:D:72:ASN:H	2.13	0.46
1:C:81:VAL:HG13	1:C:85:TRP:CB	2.46	0.46
2:M:3:DG:H2''	2:M:4:DG:C8	2.51	0.46
2:O:2:DG:C2'	2:O:3:DG:OP2	2.64	0.46
1:B:169:ARG:NH2	1:E:172:GLN:HE22	2.06	0.46
1:D:28:PHE:O	1:D:32:THR:HG23	2.16	0.46
1:E:208:PHE:C	1:E:210:LYS:N	2.74	0.46
1:H:48:VAL:HG22	1:H:138:ILE:HG23	1.97	0.46
2:M:4:DG:H2''	2:M:5:DA:C8	2.51	0.46
1:D:139:SER:HB2	1:F:85:TRP:HH2	1.81	0.45
1:A:164:THR:HG23	1:A:186:TRP:CB	2.45	0.45
1:B:86:ARG:HD3	1:B:86:ARG:HA	1.61	0.45
1:D:172:GLN:HE22	1:F:169:ARG:NH2	2.12	0.45
1:H:162:ASN:HD22	3:L:12:DC:H5	1.61	0.45
1:G:67:ASN:ND2	1:G:72:ASN:H	2.15	0.45
3:J:2:DC:H2''	3:J:3:DG:C8	2.51	0.45
1:A:140:THR:HG21	1:H:140:THR:HG21	1.91	0.45
1:B:85:TRP:CE3	1:B:86:ARG:N	2.85	0.45
1:H:62:GLU:OE2	1:H:65:ARG:NH2	2.49	0.45
1:A:19:LEU:HD23	1:H:137:MET:HE2	1.99	0.45
1:D:62:GLU:H	1:D:62:GLU:HG2	1.40	0.45
1:E:86:ARG:N	2:O:15:DG:OP1	2.49	0.45
1:H:163:HIS:ND1	3:L:13:DG:C8	2.85	0.45
1:D:16:ASP:N	1:D:16:ASP:OD1	2.49	0.45
1:D:165:PRO:CD	1:D:186:TRP:CE3	3.00	0.45
1:D:12:VAL:CG1	1:F:131:GLU:HG3	2.46	0.45
1:E:22:ILE:CD1	1:E:23:LYS:N	2.77	0.45
1:F:162:ASN:N	1:F:162:ASN:HD22	2.15	0.45
1:B:155:THR:HG22	1:E:86:ARG:O	2.16	0.45
1:H:178:ASP:OD1	1:H:184:ARG:HD2	2.17	0.45
1:A:8:ARG:HE	1:A:8:ARG:HB3	1.47	0.45
2:O:14:DG:H1'	2:O:15:DG:H5'	1.98	0.45
1:F:180:ILE:HD12	1:F:181:HIS:NE2	2.31	0.44
1:H:174:LYS:HG3	1:H:187:ALA:HB1	1.98	0.44
2:M:2:DG:H2''	2:M:3:DG:C8	2.52	0.44
1:G:202:GLU:OE2	1:G:205:LYS:NZ	2.49	0.44
1:A:2:HIS:HB3	1:H:35:HIS:HA	1.99	0.44
1:A:19:LEU:CD2	1:H:137:MET:HE2	2.47	0.44
1:C:133:THR:O	1:G:18:PHE:CE1	2.70	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:GLU:HG3	1:A:65:ARG:NH2	2.32	0.44
1:D:104:ARG:HH12	3:J:14:DG:P	2.40	0.44
1:A:165:PRO:HD3	1:A:186:TRP:CZ3	2.53	0.44
1:C:165:PRO:HD3	1:C:186:TRP:CE3	2.52	0.44
1:H:15:ASN:C	1:H:15:ASN:HD22	2.26	0.44
1:F:37:LEU:HD12	1:F:38:PRO:HA	1.99	0.44
1:B:18:PHE:CD1	1:B:22:ILE:HD11	2.52	0.44
1:A:86:ARG:CB	1:H:155:THR:O	2.67	0.43
1:D:196:GLU:HG2	4:D:403:HOH:O	2.17	0.43
1:E:60:TYR:HB3	1:E:63:TYR:HD2	1.83	0.43
1:A:24:ASP:OD2	1:H:13:TYR:OH	2.25	0.43
3:N:16:DC:H1'	3:N:17:DC:H5'	1.99	0.43
1:E:72:ASN:HB3	1:F:198:TYR:OH	2.18	0.43
1:H:198:TYR:CD2	1:H:198:TYR:C	2.95	0.43
1:E:196:GLU:OE2	1:E:196:GLU:CA	2.66	0.43
1:H:62:GLU:HG2	1:H:208:PHE:CD1	2.53	0.43
3:J:1:DG:H1'	3:J:2:DC:C6	2.54	0.43
3:P:21:DC:C2'	3:P:22:DG:O5'	2.65	0.43
1:B:140:THR:HG23	1:E:136:ASP:O	2.19	0.43
1:C:35:HIS:HD2	1:G:2:HIS:CE1	2.36	0.43
1:C:81:VAL:CG1	1:C:85:TRP:CB	2.97	0.43
1:F:180:ILE:CD1	1:F:181:HIS:NE2	2.81	0.43
1:G:169:ARG:O	1:G:172:GLN:HG2	2.18	0.43
1:C:131:GLU:O	1:C:132:ALA:C	2.61	0.43
1:E:67:ASN:HA	1:E:70:ALA:O	2.19	0.43
1:B:46:ALA:HB1	1:B:80:ALA:O	2.19	0.43
1:B:67:ASN:HD21	1:B:73:LEU:N	1.92	0.43
1:E:208:PHE:C	1:E:210:LYS:H	2.27	0.43
1:B:152:LEU:HD12	1:B:156:VAL:HB	2.01	0.43
1:F:176:ASP:HA	1:F:179:VAL:HG13	2.00	0.43
1:B:195:SER:HB2	1:B:200:ILE:HD11	2.00	0.42
1:A:19:LEU:CD2	1:H:137:MET:CE	2.96	0.42
1:H:182:PRO:HA	1:H:183:GLY:HA2	1.79	0.42
1:C:12:VAL:HG13	1:G:131:GLU:CG	2.48	0.42
1:D:96:ARG:H	1:D:97:ALA:CA	2.31	0.42
1:E:89:ARG:O	1:E:91:SER:N	2.52	0.42
1:H:104:ARG:HD2	1:H:104:ARG:HA	1.78	0.42
1:B:85:TRP:CG	1:B:86:ARG:N	2.88	0.42
1:E:86:ARG:HA	1:E:86:ARG:NE	2.34	0.42
3:J:7:DG:H2''	3:J:8:DC:OP2	2.20	0.42
3:L:17:DC:H1'	3:L:18:DT:C6	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:ASP:HA	1:A:179:VAL:CG1	2.49	0.42
1:F:53:TYR:O	1:F:63:TYR:HB3	2.18	0.42
1:D:47:GLY:HA3	1:D:128:VAL:O	2.20	0.42
1:A:140:THR:CG2	1:H:140:THR:HB	2.49	0.42
2:I:9:DC:C2'	2:I:10:DC:H5'	2.50	0.42
1:A:86:ARG:HA	1:H:155:THR:HB	2.01	0.42
1:C:52:TYR:O	1:C:123:PHE:HA	2.19	0.42
1:C:182:PRO:HA	1:C:183:GLY:HA2	1.85	0.42
1:F:210:LYS:O	1:F:211:SER:CB	2.68	0.42
1:C:169:ARG:HE	1:G:172:GLN:NE2	2.18	0.41
1:D:196:GLU:HG2	1:D:196:GLU:H	1.73	0.41
3:P:12:DC:H2''	3:P:13:DG:O5'	2.21	0.41
1:B:176:ASP:HA	1:B:179:VAL:HG13	2.03	0.41
1:D:165:PRO:HD3	1:D:186:TRP:CE3	2.55	0.41
1:E:4:LYS:HB2	1:E:5:LYS:H	1.45	0.41
1:E:51:LEU:HB2	1:E:76:TYR:HB3	2.02	0.41
1:F:16:ASP:C	1:F:18:PHE:H	2.28	0.41
1:F:79:LYS:HD3	1:F:142:GLN:OE1	2.20	0.41
1:E:86:ARG:NH2	4:E:361:HOH:O	2.53	0.41
1:G:18:PHE:HE2	1:G:140:THR:HG21	1.86	0.41
1:C:44:GLN:HG2	1:C:97:ALA:O	2.20	0.41
1:F:165:PRO:HD3	1:F:186:TRP:CE3	2.55	0.41
1:D:50:ALA:O	1:D:125:CYS:HA	2.20	0.41
1:E:208:PHE:O	1:E:210:LYS:N	2.53	0.41
1:A:150:LYS:N	1:A:151:PRO:HD3	2.36	0.41
1:A:165:PRO:HD3	1:A:186:TRP:CE3	2.55	0.41
1:B:36:SER:OG	1:E:2:HIS:CE1	2.74	0.41
1:F:18:PHE:CE2	1:F:19:LEU:HD13	2.56	0.41
3:N:7:DG:H2''	3:N:8:DC:OP2	2.21	0.41
1:C:168:GLY:HA2	1:G:171:ALA:C	2.46	0.41
1:G:42:ARG:HD3	1:G:106:ARG:NE	2.36	0.41
2:K:18:DG:H1	3:L:5:DC:H42	1.68	0.41
1:A:46:ALA:HB1	1:A:80:ALA:O	2.20	0.40
1:D:119:ASP:HB2	4:D:221:HOH:O	2.20	0.40
1:F:14:ARG:HE	1:F:14:ARG:HB2	1.67	0.40
1:H:172:GLN:O	1:H:190:CYS:HB3	2.22	0.40
1:D:100:GLU:H	1:D:100:GLU:CD	2.30	0.40
1:G:73:LEU:HA	1:G:74:PRO:HD2	1.88	0.40
3:N:3:DG:H2''	3:N:4:DG:C8	2.56	0.40
1:B:66:ILE:HG21	1:B:66:ILE:HD13	1.81	0.40
1:C:172:GLN:NE2	1:G:169:ARG:NH2	2.63	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:6:PHE:CD2	1:H:6:PHE:C	3.00	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:1:DC:OP1	2:O:21:DG:OP2[1_556]	1.97	0.23
1:A:133:THR:OG1	1:G:61:ASP:OD1[2_656]	2.05	0.15

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	211/235 (90%)	194 (92%)	11 (5%)	6 (3%)	4 6
1	B	211/235 (90%)	197 (93%)	8 (4%)	6 (3%)	4 6
1	C	206/235 (88%)	189 (92%)	12 (6%)	5 (2%)	4 8
1	D	206/235 (88%)	188 (91%)	13 (6%)	5 (2%)	4 8
1	E	209/235 (89%)	189 (90%)	13 (6%)	7 (3%)	3 4
1	F	208/235 (88%)	192 (92%)	12 (6%)	4 (2%)	6 11
1	G	206/235 (88%)	191 (93%)	12 (6%)	3 (2%)	8 16
1	H	207/235 (88%)	187 (90%)	17 (8%)	3 (1%)	9 17
All	All	1664/1880 (88%)	1527 (92%)	98 (6%)	39 (2%)	5 8

All (39) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	89	ARG
1	A	93	HIS
1	B	4	LYS

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Mol	Chain	Res	Type
1	B	95	THR
1	C	88	SER
1	C	90	ILE
1	E	88	SER
1	E	90	ILE
1	F	94	GLU
1	F	96	ARG
1	H	88	SER
1	H	89	ARG
1	A	84	GLY
1	A	95	THR
1	B	84	GLY
1	B	91	SER
1	B	92	ASP
1	C	93	HIS
1	C	132	ALA
1	D	133	THR
1	E	89	ARG
1	E	95	THR
1	E	98	GLY
1	E	209	SER
1	G	3	ASN
1	C	5	LYS
1	D	10	GLU
1	F	17	SER
1	H	87	GLN
1	B	93	HIS
1	D	89	ARG
1	D	93	HIS
1	E	2	HIS
1	F	5	LYS
1	G	91	SER
1	A	86	ARG
1	D	132	ALA
1	G	29	PHE
1	A	90	ILE

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	170/201 (85%)	154 (91%)	16 (9%)	8	18
1	B	170/201 (85%)	155 (91%)	15 (9%)	9	20
1	C	167/201 (83%)	151 (90%)	16 (10%)	8	17
1	D	165/201 (82%)	151 (92%)	14 (8%)	10	22
1	E	164/201 (82%)	147 (90%)	17 (10%)	7	14
1	F	172/201 (86%)	153 (89%)	19 (11%)	6	13
1	G	171/201 (85%)	149 (87%)	22 (13%)	4	9
1	H	168/201 (84%)	150 (89%)	18 (11%)	6	13
All	All	1347/1608 (84%)	1210 (90%)	137 (10%)	7	15

All (137) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	HIS
1	A	8	ARG
1	A	14	ARG
1	A	16	ASP
1	A	42	ARG
1	A	51	LEU
1	A	81	VAL
1	A	89	ARG
1	A	140	THR
1	A	164	THR
1	A	174	LYS
1	A	179	VAL
1	A	185	GLU
1	A	188	GLU
1	A	201	GLU
1	A	205	LYS
1	B	2	HIS
1	B	5	LYS
1	B	20	GLU
1	B	21	LEU
1	B	27	ARG
1	B	41	GLU
1	B	42	ARG
1	B	51	LEU

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Mol	Chain	Res	Type
1	B	81	VAL
1	B	85	TRP
1	B	100	GLU
1	B	110	ARG
1	B	179	VAL
1	B	185	GLU
1	B	196	GLU
1	C	5	LYS
1	C	8	ARG
1	C	14	ARG
1	C	16	ASP
1	C	19	LEU
1	C	21	LEU
1	C	51	LEU
1	C	62	GLU
1	C	65	ARG
1	C	81	VAL
1	C	100	GLU
1	C	104	ARG
1	C	179	VAL
1	C	180	ILE
1	C	201	GLU
1	C	206	GLN
1	D	8	ARG
1	D	16	ASP
1	D	19	LEU
1	D	21	LEU
1	D	51	LEU
1	D	62	GLU
1	D	81	VAL
1	D	100	GLU
1	D	135	SER
1	D	179	VAL
1	D	180	ILE
1	D	185	GLU
1	D	196	GLU
1	D	210	LYS
1	E	2	HIS
1	E	4	LYS
1	E	15	ASN
1	E	21	LEU
1	E	22	ILE

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Mol	Chain	Res	Type
1	E	27	ARG
1	E	42	ARG
1	E	51	LEU
1	E	81	VAL
1	E	93	HIS
1	E	100	GLU
1	E	104	ARG
1	E	124	SER
1	E	133	THR
1	E	135	SER
1	E	179	VAL
1	E	188	GLU
1	F	15	ASN
1	F	18	PHE
1	F	19	LEU
1	F	20	GLU
1	F	21	LEU
1	F	27	ARG
1	F	42	ARG
1	F	51	LEU
1	F	81	VAL
1	F	85	TRP
1	F	90	ILE
1	F	100	GLU
1	F	106	ARG
1	F	133	THR
1	F	164	THR
1	F	179	VAL
1	F	184	ARG
1	F	205	LYS
1	F	210	LYS
1	G	2	HIS
1	G	3	ASN
1	G	4	LYS
1	G	15	ASN
1	G	18	PHE
1	G	21	LEU
1	G	51	LEU
1	G	68	ARG
1	G	85	TRP
1	G	90	ILE
1	G	91	SER

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Mol	Chain	Res	Type
1	G	94	GLU
1	G	95	THR
1	G	99	SER
1	G	100	GLU
1	G	114	LYS
1	G	139	SER
1	G	174	LYS
1	G	179	VAL
1	G	201	GLU
1	G	205	LYS
1	G	206	GLN
1	H	2	HIS
1	H	15	ASN
1	H	21	LEU
1	H	27	ARG
1	H	51	LEU
1	H	85	TRP
1	H	93	HIS
1	H	100	GLU
1	H	104	ARG
1	H	133	THR
1	H	140	THR
1	H	172	GLN
1	H	179	VAL
1	H	180	ILE
1	H	188	GLU
1	H	203	ARG
1	H	209	SER
1	H	210	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (61) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	HIS
1	A	35	HIS
1	A	44	GLN
1	A	67	ASN
1	A	103	ASN
1	A	162	ASN
1	A	172	GLN
1	A	206	GLN
1	B	2	HIS

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Mol	Chain	Res	Type
1	B	35	HIS
1	B	44	GLN
1	B	67	ASN
1	B	93	HIS
1	B	103	ASN
1	B	142	GLN
1	B	162	ASN
1	B	172	GLN
1	C	35	HIS
1	C	44	GLN
1	C	67	ASN
1	C	103	ASN
1	C	108	HIS
1	C	162	ASN
1	C	172	GLN
1	D	44	GLN
1	D	67	ASN
1	D	103	ASN
1	D	108	HIS
1	D	162	ASN
1	D	172	GLN
1	D	206	GLN
1	E	2	HIS
1	E	15	ASN
1	E	35	HIS
1	E	44	GLN
1	E	67	ASN
1	E	103	ASN
1	E	111	ASN
1	E	162	ASN
1	E	172	GLN
1	F	15	ASN
1	F	35	HIS
1	F	44	GLN
1	F	67	ASN
1	F	103	ASN
1	F	162	ASN
1	F	172	GLN
1	G	15	ASN
1	G	67	ASN
1	G	87	GLN
1	G	103	ASN

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Mol	Chain	Res	Type
1	G	172	GLN
1	H	2	HIS
1	H	15	ASN
1	H	35	HIS
1	H	44	GLN
1	H	67	ASN
1	H	103	ASN
1	H	111	ASN
1	H	162	ASN
1	H	172	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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 [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	213/235 (90%)	0.02	10 (4%) 36 32	27, 42, 96, 111	0
1	B	213/235 (90%)	-0.06	11 (5%) 33 29	28, 42, 92, 109	0
1	C	208/235 (88%)	0.12	7 (3%) 48 43	35, 56, 92, 118	0
1	D	208/235 (88%)	0.17	6 (2%) 53 49	34, 54, 91, 111	0
1	E	211/235 (89%)	0.00	9 (4%) 40 35	29, 46, 97, 130	0
1	F	210/235 (89%)	-0.02	6 (2%) 53 49	30, 48, 86, 107	0
1	G	208/235 (88%)	-0.10	4 (1%) 66 62	27, 44, 85, 108	0
1	H	209/235 (88%)	0.16	12 (5%) 29 26	33, 52, 98, 133	0
2	I	22/22 (100%)	0.43	3 (13%) 7 5	31, 74, 142, 183	0
2	K	22/22 (100%)	0.51	0 100 100	30, 73, 133, 155	0
2	M	22/22 (100%)	0.23	1 (4%) 38 33	32, 65, 124, 142	0
2	O	22/22 (100%)	0.26	0 100 100	29, 66, 118, 126	0
3	J	22/22 (100%)	0.38	1 (4%) 38 33	33, 77, 129, 149	0
3	L	22/22 (100%)	0.59	3 (13%) 7 5	30, 80, 144, 156	0
3	N	22/22 (100%)	0.25	2 (9%) 15 13	28, 72, 130, 172	0
3	P	22/22 (100%)	0.33	0 100 100	27, 71, 119, 132	0
All	All	1856/2056 (90%)	0.07	75 (4%) 42 38	27, 49, 104, 183	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	214	THR	6.2
1	H	90	ILE	5.1
1	A	93	HIS	4.7
1	B	91	SER	4.7
1	B	214	THR	4.7

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Mol	Chain	Res	Type	RSRZ
1	E	211	SER	4.4
1	H	85	TRP	4.1
2	M	1	DC	3.8
1	E	85	TRP	3.8
1	E	92	ASP	3.8
1	D	89	ARG	3.7
1	C	95	THR	3.6
3	N	1	DG	3.5
1	D	6	PHE	3.4
1	B	87	GLN	3.4
1	E	90	ILE	3.4
1	B	3	ASN	3.4
1	H	91	SER	3.2
1	F	92	ASP	3.1
1	C	211	SER	3.1
1	H	210	LYS	3.0
1	C	201	GLU	3.0
1	D	212	ASN	3.0
1	A	94	GLU	2.9
1	A	90	ILE	2.9
1	B	90	ILE	2.9
1	B	94	GLU	2.9
1	G	93	HIS	2.8
1	B	93	HIS	2.8
1	H	135	SER	2.8
1	C	97	ALA	2.8
2	I	2	DG	2.7
3	J	1	DG	2.7
1	H	27	ARG	2.6
1	E	95	THR	2.6
1	G	97	ALA	2.6
1	G	92	ASP	2.5
1	H	87	GLN	2.5
1	E	3	ASN	2.5
1	A	91	SER	2.5
1	D	88	SER	2.5
1	C	85	TRP	2.5
1	E	93	HIS	2.5
1	A	89	ARG	2.5
1	F	97	ALA	2.5
1	B	88	SER	2.5
1	D	90	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
3	N	2	DC	2.4
1	A	3	ASN	2.4
1	F	211	SER	2.4
1	E	209	SER	2.4
1	F	2	HIS	2.3
1	B	92	ASP	2.3
2	I	1	DC	2.3
1	H	5	LYS	2.3
1	F	4	LYS	2.3
1	F	96	ARG	2.3
3	L	1	DG	2.2
1	H	58	SER	2.2
1	C	93	HIS	2.2
1	H	2	HIS	2.2
1	A	92	ASP	2.2
1	D	85	TRP	2.2
1	H	209	SER	2.2
1	H	95	THR	2.2
1	A	85	TRP	2.1
3	L	22	DG	2.1
1	B	89	ARG	2.1
1	C	65	ARG	2.1
1	G	2	HIS	2.1
1	B	96	ARG	2.1
3	L	2	DC	2.1
2	I	5	DA	2.0
1	E	1	GLY	2.0
1	A	95	THR	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.