



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:05 PM GMT

PDB ID : 4NOD
Title : Distinct structural features of TFAM drive mitochondrial DNA packaging versus transcriptional activation
Authors : Ngo, H.B.; Lovely, G.A.; Phillips, R.; Chan, D.C.
Deposited on : 2013-11-19
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

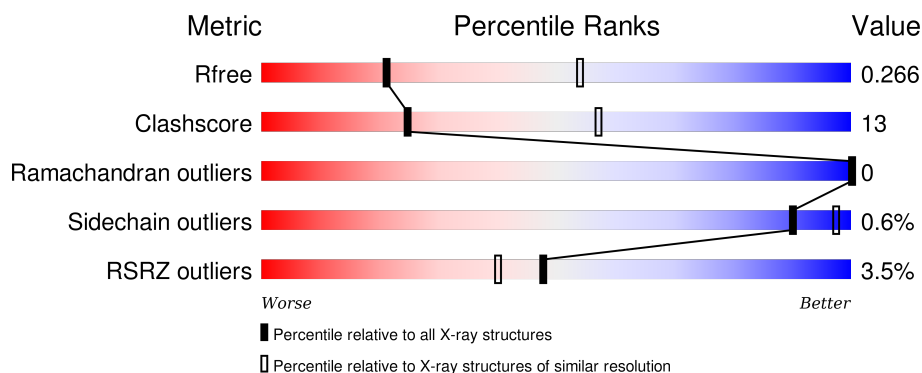
1 Overall quality at a glance

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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	230	<div> <div>7%</div> <div>67% 14% • 17%</div> </div>
1	B	230	<div> <div>6%</div> <div>62% 19% • 17%</div> </div>
1	G	230	<div> <div>66% 15% • 17%</div> </div>
1	H	230	<div> <div>63% 18% • 17%</div> </div>
2	C	22	<div> <div>50% 50%</div> </div>

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Mol	Chain	Length	Quality of chain
2	E	22	<div><div></div><div>50%</div><div></div><div>50%</div><div></div></div>
2	I	22	<div><div></div><div>50%</div><div></div><div>50%</div><div></div></div>
2	K	22	<div><div></div><div>55%</div><div></div><div>45%</div><div></div></div>
3	D	22	<div><div></div><div>32%</div><div></div><div>64%</div><div></div><div>5%</div></div>
3	F	22	<div><div>5%</div><div></div><div>27%</div><div></div><div>68%</div><div></div><div>5%</div></div>
3	J	22	<div><div></div><div>36%</div><div></div><div>59%</div><div></div><div>5%</div></div>
3	L	22	<div><div></div><div>41%</div><div></div><div>55%</div><div></div><div>5%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9985 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 Transcription factor A, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	190	Total	C	N	O	S	0	0	0
			1602	1012	289	295	6			
1	B	190	Total	C	N	O	S	0	0	0
			1602	1012	289	295	6			
1	G	190	Total	C	N	O	S	0	0	0
			1602	1012	289	295	6			
1	H	190	Total	C	N	O	S	0	0	0
			1602	1012	289	295	6			

There are 140 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	8	MET	-	EXPRESSION TAG	UNP Q00059
A	9	GLY	-	EXPRESSION TAG	UNP Q00059
A	10	SER	-	EXPRESSION TAG	UNP Q00059
A	11	SER	-	EXPRESSION TAG	UNP Q00059
A	12	HIS	-	EXPRESSION TAG	UNP Q00059
A	13	HIS	-	EXPRESSION TAG	UNP Q00059
A	14	HIS	-	EXPRESSION TAG	UNP Q00059
A	15	HIS	-	EXPRESSION TAG	UNP Q00059
A	16	HIS	-	EXPRESSION TAG	UNP Q00059
A	17	HIS	-	EXPRESSION TAG	UNP Q00059
A	18	SER	-	EXPRESSION TAG	UNP Q00059
A	19	SER	-	EXPRESSION TAG	UNP Q00059
A	20	GLY	-	EXPRESSION TAG	UNP Q00059
A	21	LEU	-	EXPRESSION TAG	UNP Q00059
A	22	VAL	-	EXPRESSION TAG	UNP Q00059
A	23	PRO	-	EXPRESSION TAG	UNP Q00059
A	24	ARG	-	EXPRESSION TAG	UNP Q00059
A	25	GLY	-	EXPRESSION TAG	UNP Q00059
A	26	SER	-	EXPRESSION TAG	UNP Q00059
A	27	HIS	-	EXPRESSION TAG	UNP Q00059
A	28	MET	-	EXPRESSION TAG	UNP Q00059

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Chain	Residue	Modelled	Actual	Comment	Reference
A	29	ALA	-	EXPRESSION TAG	UNP Q00059
A	30	SER	-	EXPRESSION TAG	UNP Q00059
A	31	MET	-	EXPRESSION TAG	UNP Q00059
A	32	THR	-	EXPRESSION TAG	UNP Q00059
A	33	GLY	-	EXPRESSION TAG	UNP Q00059
A	34	GLY	-	EXPRESSION TAG	UNP Q00059
A	35	GLN	-	EXPRESSION TAG	UNP Q00059
A	36	GLN	-	EXPRESSION TAG	UNP Q00059
A	37	MET	-	EXPRESSION TAG	UNP Q00059
A	38	GLY	-	EXPRESSION TAG	UNP Q00059
A	39	ARG	-	EXPRESSION TAG	UNP Q00059
A	40	GLY	-	EXPRESSION TAG	UNP Q00059
A	41	SER	-	EXPRESSION TAG	UNP Q00059
A	42	MET	-	EXPRESSION TAG	UNP Q00059
B	8	MET	-	EXPRESSION TAG	UNP Q00059
B	9	GLY	-	EXPRESSION TAG	UNP Q00059
B	10	SER	-	EXPRESSION TAG	UNP Q00059
B	11	SER	-	EXPRESSION TAG	UNP Q00059
B	12	HIS	-	EXPRESSION TAG	UNP Q00059
B	13	HIS	-	EXPRESSION TAG	UNP Q00059
B	14	HIS	-	EXPRESSION TAG	UNP Q00059
B	15	HIS	-	EXPRESSION TAG	UNP Q00059
B	16	HIS	-	EXPRESSION TAG	UNP Q00059
B	17	HIS	-	EXPRESSION TAG	UNP Q00059
B	18	SER	-	EXPRESSION TAG	UNP Q00059
B	19	SER	-	EXPRESSION TAG	UNP Q00059
B	20	GLY	-	EXPRESSION TAG	UNP Q00059
B	21	LEU	-	EXPRESSION TAG	UNP Q00059
B	22	VAL	-	EXPRESSION TAG	UNP Q00059
B	23	PRO	-	EXPRESSION TAG	UNP Q00059
B	24	ARG	-	EXPRESSION TAG	UNP Q00059
B	25	GLY	-	EXPRESSION TAG	UNP Q00059
B	26	SER	-	EXPRESSION TAG	UNP Q00059
B	27	HIS	-	EXPRESSION TAG	UNP Q00059
B	28	MET	-	EXPRESSION TAG	UNP Q00059
B	29	ALA	-	EXPRESSION TAG	UNP Q00059
B	30	SER	-	EXPRESSION TAG	UNP Q00059
B	31	MET	-	EXPRESSION TAG	UNP Q00059
B	32	THR	-	EXPRESSION TAG	UNP Q00059
B	33	GLY	-	EXPRESSION TAG	UNP Q00059
B	34	GLY	-	EXPRESSION TAG	UNP Q00059
B	35	GLN	-	EXPRESSION TAG	UNP Q00059

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Chain	Residue	Modelled	Actual	Comment	Reference
B	36	GLN	-	EXPRESSION TAG	UNP Q00059
B	37	MET	-	EXPRESSION TAG	UNP Q00059
B	38	GLY	-	EXPRESSION TAG	UNP Q00059
B	39	ARG	-	EXPRESSION TAG	UNP Q00059
B	40	GLY	-	EXPRESSION TAG	UNP Q00059
B	41	SER	-	EXPRESSION TAG	UNP Q00059
B	42	MET	-	EXPRESSION TAG	UNP Q00059
G	8	MET	-	EXPRESSION TAG	UNP Q00059
G	9	GLY	-	EXPRESSION TAG	UNP Q00059
G	10	SER	-	EXPRESSION TAG	UNP Q00059
G	11	SER	-	EXPRESSION TAG	UNP Q00059
G	12	HIS	-	EXPRESSION TAG	UNP Q00059
G	13	HIS	-	EXPRESSION TAG	UNP Q00059
G	14	HIS	-	EXPRESSION TAG	UNP Q00059
G	15	HIS	-	EXPRESSION TAG	UNP Q00059
G	16	HIS	-	EXPRESSION TAG	UNP Q00059
G	17	HIS	-	EXPRESSION TAG	UNP Q00059
G	18	SER	-	EXPRESSION TAG	UNP Q00059
G	19	SER	-	EXPRESSION TAG	UNP Q00059
G	20	GLY	-	EXPRESSION TAG	UNP Q00059
G	21	LEU	-	EXPRESSION TAG	UNP Q00059
G	22	VAL	-	EXPRESSION TAG	UNP Q00059
G	23	PRO	-	EXPRESSION TAG	UNP Q00059
G	24	ARG	-	EXPRESSION TAG	UNP Q00059
G	25	GLY	-	EXPRESSION TAG	UNP Q00059
G	26	SER	-	EXPRESSION TAG	UNP Q00059
G	27	HIS	-	EXPRESSION TAG	UNP Q00059
G	28	MET	-	EXPRESSION TAG	UNP Q00059
G	29	ALA	-	EXPRESSION TAG	UNP Q00059
G	30	SER	-	EXPRESSION TAG	UNP Q00059
G	31	MET	-	EXPRESSION TAG	UNP Q00059
G	32	THR	-	EXPRESSION TAG	UNP Q00059
G	33	GLY	-	EXPRESSION TAG	UNP Q00059
G	34	GLY	-	EXPRESSION TAG	UNP Q00059
G	35	GLN	-	EXPRESSION TAG	UNP Q00059
G	36	GLN	-	EXPRESSION TAG	UNP Q00059
G	37	MET	-	EXPRESSION TAG	UNP Q00059
G	38	GLY	-	EXPRESSION TAG	UNP Q00059
G	39	ARG	-	EXPRESSION TAG	UNP Q00059
G	40	GLY	-	EXPRESSION TAG	UNP Q00059
G	41	SER	-	EXPRESSION TAG	UNP Q00059
G	42	MET	-	EXPRESSION TAG	UNP Q00059

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Chain	Residue	Modelled	Actual	Comment	Reference
H	8	MET	-	EXPRESSION TAG	UNP Q00059
H	9	GLY	-	EXPRESSION TAG	UNP Q00059
H	10	SER	-	EXPRESSION TAG	UNP Q00059
H	11	SER	-	EXPRESSION TAG	UNP Q00059
H	12	HIS	-	EXPRESSION TAG	UNP Q00059
H	13	HIS	-	EXPRESSION TAG	UNP Q00059
H	14	HIS	-	EXPRESSION TAG	UNP Q00059
H	15	HIS	-	EXPRESSION TAG	UNP Q00059
H	16	HIS	-	EXPRESSION TAG	UNP Q00059
H	17	HIS	-	EXPRESSION TAG	UNP Q00059
H	18	SER	-	EXPRESSION TAG	UNP Q00059
H	19	SER	-	EXPRESSION TAG	UNP Q00059
H	20	GLY	-	EXPRESSION TAG	UNP Q00059
H	21	LEU	-	EXPRESSION TAG	UNP Q00059
H	22	VAL	-	EXPRESSION TAG	UNP Q00059
H	23	PRO	-	EXPRESSION TAG	UNP Q00059
H	24	ARG	-	EXPRESSION TAG	UNP Q00059
H	25	GLY	-	EXPRESSION TAG	UNP Q00059
H	26	SER	-	EXPRESSION TAG	UNP Q00059
H	27	HIS	-	EXPRESSION TAG	UNP Q00059
H	28	MET	-	EXPRESSION TAG	UNP Q00059
H	29	ALA	-	EXPRESSION TAG	UNP Q00059
H	30	SER	-	EXPRESSION TAG	UNP Q00059
H	31	MET	-	EXPRESSION TAG	UNP Q00059
H	32	THR	-	EXPRESSION TAG	UNP Q00059
H	33	GLY	-	EXPRESSION TAG	UNP Q00059
H	34	GLY	-	EXPRESSION TAG	UNP Q00059
H	35	GLN	-	EXPRESSION TAG	UNP Q00059
H	36	GLN	-	EXPRESSION TAG	UNP Q00059
H	37	MET	-	EXPRESSION TAG	UNP Q00059
H	38	GLY	-	EXPRESSION TAG	UNP Q00059
H	39	ARG	-	EXPRESSION TAG	UNP Q00059
H	40	GLY	-	EXPRESSION TAG	UNP Q00059
H	41	SER	-	EXPRESSION TAG	UNP Q00059
H	42	MET	-	EXPRESSION TAG	UNP Q00059

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	22	Total	Br	C	N	O	P	0	0	0
			460	1	218	84	136	21			

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	E	22	Total	Br	C	N	O	P	0	0	0
			460	1	218	84	136	21			
2	I	22	Total	Br	C	N	O	P	0	0	0
			460	1	218	84	136	21			
2	K	22	Total	Br	C	N	O	P	0	0	0
			459	1	217	84	136	21			

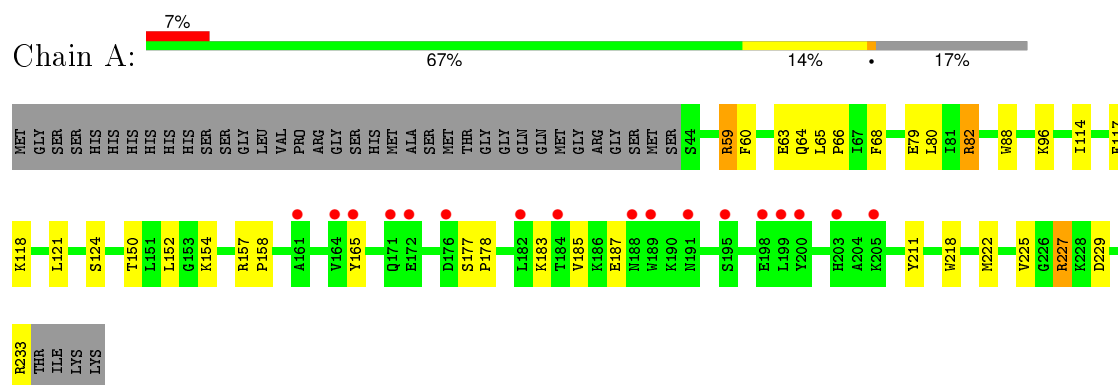
- Molecule 3 is a DNA chain called 5'-D(*CP*CP*AP*AP*CP*CP*AP*AP*GP*CP*CP*CP*CP*AP*TP*AP*CP*CP*CP*CP*AP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	22	Total	C	N	O	P	0	0	0
			435	208	83	123	21			
3	F	22	Total	C	N	O	P	0	0	0
			435	208	83	123	21			
3	J	22	Total	C	N	O	P	0	0	0
			434	208	83	122	21			
3	L	22	Total	C	N	O	P	0	0	0
			434	208	83	122	21			

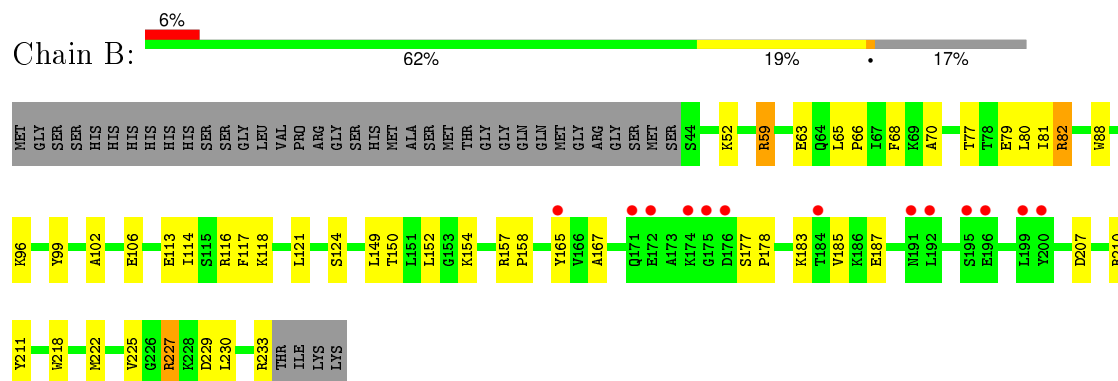
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

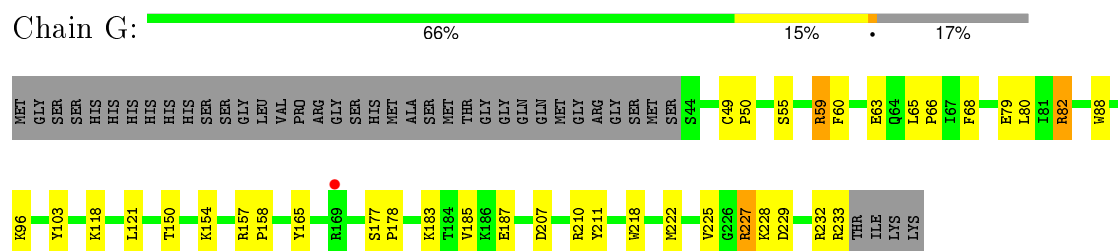
- Molecule 1: Transcription factor A, mitochondrial



- Molecule 1: Transcription factor A, mitochondrial

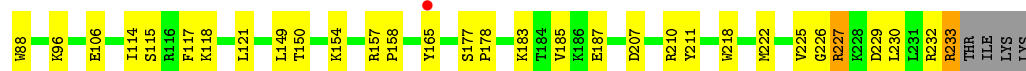
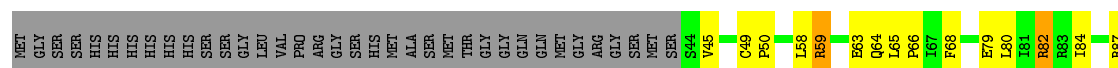


- Molecule 1: Transcription factor A, mitochondrial



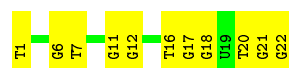
- Molecule 1: Transcription factor A, mitochondrial

Chain H: 



• Molecule 2: 5'-D(*TP*TP*GP*GP*GP*GP*TP*AP*TP*GP*GP*GP*GP*CP*TP*TP*GP*GP*(BRU)P*TP*GP*G)-3'

Chain C: 



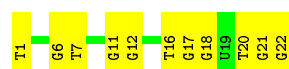
• Molecule 2: 5'-D(*TP*TP*GP*GP*GP*GP*TP*AP*TP*GP*GP*GP*GP*CP*TP*TP*GP*GP*(BRU)P*TP*GP*G)-3'

Chain E: 



• Molecule 2: 5'-D(*TP*TP*GP*GP*GP*GP*TP*AP*TP*GP*GP*GP*GP*CP*TP*TP*GP*GP*(BRU)P*TP*GP*G)-3'

Chain I: 



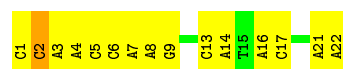
• Molecule 2: 5'-D(*TP*TP*GP*GP*GP*GP*TP*AP*TP*GP*GP*GP*GP*CP*TP*TP*GP*GP*(BRU)P*TP*GP*G)-3'

Chain K: 

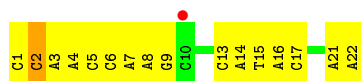


• Molecule 3: 5'-D(*CP*CP*AP*AP*CP*CP*AP*AP*GP*CP*CP*CP*CP*AP*TP*AP*CP*C P*CP*CP*AP*A)-3'

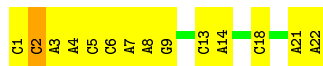
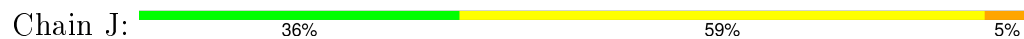
Chain D: 



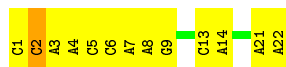
• Molecule 3: 5'-D(*CP*CP*AP*AP*CP*CP*AP*AP*GP*CP*CP*CP*CP*AP*TP*AP*CP*C P*CP*CP*AP*A)-3'



● Molecule 3: 5'-D(*CP*CP*AP*AP*CP*CP*AP*AP*GP*CP*CP*CP*CP*AP*TP*AP*CP*CP*CP*CP*AP*A)-3'



● Molecule 3: 5'-D(*CP*CP*AP*AP*CP*CP*AP*AP*GP*CP*CP*CP*CP*AP*TP*AP*CP*CP*CP*CP*AP*A)-3'



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.37Å 82.49Å 104.22Å 79.86° 85.48° 84.53°	Depositor
Resolution (Å)	35.64 – 2.90 35.62 – 2.90	Depositor EDS
% Data completeness (in resolution range)	87.0 (35.64-2.90) 82.4 (35.62-2.90)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.221 , 0.267 0.220 , 0.266	Depositor DCC
R_{free} test set	1783 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	79.1	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 50.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 35526 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9985	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.60% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BRU

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	0.52	0/1633	0.83	7/2184 (0.3%)
1	B	0.58	0/1633	1.03	11/2184 (0.5%)
1	G	0.62	0/1633	0.90	10/2184 (0.5%)
1	H	0.58	0/1633	0.88	8/2184 (0.4%)
2	C	0.33	0/493	0.76	0/762
2	E	0.35	0/493	0.77	0/762
2	I	0.43	0/493	0.79	0/762
2	K	0.40	0/490	0.76	0/755
3	D	0.46	0/487	0.91	2/745 (0.3%)
3	F	0.45	0/487	0.91	2/745 (0.3%)
3	J	0.55	1/486 (0.2%)	0.92	2/744 (0.3%)
3	L	0.51	0/486	0.90	2/744 (0.3%)
All	All	0.53	1/10447 (0.0%)	0.89	44/14755 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	J	18	DC	O3'-P	-5.06	1.55	1.61

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	82	ARG	NE-CZ-NH2	-13.27	113.67	120.30
1	B	82	ARG	NE-CZ-NH1	12.59	126.60	120.30
1	B	227	ARG	NE-CZ-NH2	12.30	126.45	120.30
1	B	227	ARG	NE-CZ-NH1	-12.03	114.28	120.30
1	A	233	ARG	NE-CZ-NH2	11.18	125.89	120.30
1	B	233	ARG	NE-CZ-NH2	-11.00	114.80	120.30
1	A	233	ARG	NE-CZ-NH1	-10.85	114.88	120.30
1	B	233	ARG	NE-CZ-NH1	10.44	125.52	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	233	ARG	NE-CZ-NH2	-10.27	115.17	120.30
1	H	59	ARG	NE-CZ-NH1	10.10	125.35	120.30
1	G	233	ARG	NE-CZ-NH2	10.02	125.31	120.30
1	G	233	ARG	NE-CZ-NH1	-9.94	115.33	120.30
1	B	59	ARG	NE-CZ-NH1	9.73	125.17	120.30
1	H	233	ARG	NE-CZ-NH1	9.45	125.03	120.30
1	B	59	ARG	NE-CZ-NH2	-9.43	115.59	120.30
1	H	59	ARG	NE-CZ-NH2	-9.38	115.61	120.30
1	B	82	ARG	CD-NE-CZ	8.01	134.82	123.60
1	G	82	ARG	NE-CZ-NH2	7.72	124.16	120.30
1	G	59	ARG	NE-CZ-NH1	-7.46	116.57	120.30
1	A	59	ARG	NE-CZ-NH1	-7.42	116.59	120.30
1	H	82	ARG	NE-CZ-NH1	-7.29	116.66	120.30
1	G	227	ARG	NE-CZ-NH2	-7.08	116.76	120.30
1	G	59	ARG	NE-CZ-NH2	7.07	123.84	120.30
1	G	227	ARG	NE-CZ-NH1	6.96	123.78	120.30
1	G	82	ARG	NE-CZ-NH1	-6.78	116.91	120.30
1	G	232	ARG	NE-CZ-NH2	-6.78	116.91	120.30
3	F	21	DA	C4'-C3'-O3'	6.55	126.08	109.70
3	L	21	DA	C4'-C3'-O3'	6.54	126.05	109.70
1	H	82	ARG	NE-CZ-NH2	6.47	123.54	120.30
3	J	21	DA	C4'-C3'-O3'	6.47	125.86	109.70
3	D	2	DC	C1'-O4'-C4'	-6.40	103.70	110.10
1	B	227	ARG	CD-NE-CZ	6.38	132.54	123.60
3	D	21	DA	C4'-C3'-O3'	6.36	125.61	109.70
1	A	59	ARG	NE-CZ-NH2	5.94	123.27	120.30
3	L	2	DC	C1'-O4'-C4'	-5.72	104.38	110.10
1	A	227	ARG	NE-CZ-NH2	-5.71	117.44	120.30
3	J	2	DC	C1'-O4'-C4'	-5.21	104.89	110.10
1	G	233	ARG	CD-NE-CZ	5.13	130.78	123.60
1	A	82	ARG	NE-CZ-NH2	5.09	122.85	120.30
1	B	233	ARG	CD-NE-CZ	5.07	130.70	123.60
1	H	227	ARG	NE-CZ-NH1	5.07	122.83	120.30
3	F	2	DC	C1'-O4'-C4'	-5.05	105.05	110.10
1	H	227	ARG	NE-CZ-NH2	-5.04	117.78	120.30
1	A	233	ARG	CD-NE-CZ	5.01	130.61	123.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1602	0	1643	23	0
1	B	1602	0	1643	33	0
1	G	1602	0	1643	25	0
1	H	1602	0	1643	35	0
2	C	460	0	246	26	0
2	E	460	0	246	25	0
2	I	460	0	246	22	0
2	K	459	0	243	26	0
3	D	435	0	245	20	0
3	F	435	0	245	24	0
3	J	434	0	242	14	0
3	L	434	0	242	14	0
All	All	9985	0	8527	232	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 13.

All (232) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:1:DT:C5'	2:K:22:DG:H2''	1.68	1.23
2:I:22:DG:H2''	2:K:1:DT:C5'	1.75	1.17
3:J:1:DC:C5'	3:L:22:DA:O3'	2.11	0.97
3:J:22:DA:O3'	3:L:1:DC:C5'	2.13	0.97
3:D:22:DA:O3'	3:F:1:DC:C5'	2.13	0.96
2:C:1:DT:C5'	2:E:22:DG:H2''	1.98	0.94
3:D:1:DC:C5'	3:F:22:DA:O3'	2.17	0.92
2:C:22:DG:H2''	2:E:1:DT:C5'	1.99	0.91
3:D:22:DA:O3'	3:F:1:DC:H5''	1.69	0.91
3:D:1:DC:H5'	3:F:22:DA:O3'	1.73	0.89
3:D:1:DC:O5'	3:F:22:DA:O3'	1.90	0.89
2:C:1:DT:C5'	2:E:22:DG:O3'	2.21	0.88
2:I:1:DT:C5'	2:K:22:DG:C2'	2.53	0.86
1:G:79:GLU:OE1	1:G:82:ARG:NH1	2.09	0.86
1:B:79:GLU:OE1	1:B:82:ARG:NH1	2.09	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:22:DG:O3'	2:E:1:DT:C5'	2.23	0.85
3:D:22:DA:O3'	3:F:1:DC:O5'	1.96	0.84
2:C:22:DG:H1	3:D:1:DC:H42	1.24	0.83
3:D:1:DC:C5'	3:F:22:DA:HO3'	1.91	0.81
2:K:22:DG:H8	2:K:22:DG:O5'	1.63	0.80
2:I:22:DG:C2'	2:K:1:DT:C5'	2.59	0.79
3:D:22:DA:H2''	3:F:1:DC:H5''	1.64	0.79
2:I:22:DG:O5'	2:I:22:DG:H8	1.65	0.79
1:A:79:GLU:OE1	1:A:82:ARG:NH1	2.16	0.78
1:H:183:LYS:O	1:H:187:GLU:HG2	1.85	0.77
1:B:157:ARG:HB2	3:F:4:DA:H5'	1.65	0.77
2:K:2:DT:N1	2:K:2:DT:C2'	2.49	0.76
1:H:79:GLU:OE1	1:H:82:ARG:NH1	2.18	0.76
2:K:22:DG:H1	3:L:1:DC:H42	1.34	0.76
1:G:183:LYS:O	1:G:187:GLU:HG2	1.86	0.76
1:A:183:LYS:O	1:A:187:GLU:HG2	1.86	0.76
1:B:183:LYS:O	1:B:187:GLU:HG2	1.85	0.75
2:E:22:DG:H8	2:E:22:DG:O5'	1.69	0.74
2:I:18:DG:H1	3:J:5:DC:H42	1.36	0.73
2:I:22:DG:O3'	2:K:1:DT:C5'	2.38	0.71
1:G:65:LEU:HB3	1:G:66:PRO:HD3	1.72	0.70
1:G:157:ARG:HB2	3:J:4:DA:H5'	1.74	0.70
2:C:22:DG:O5'	2:C:22:DG:H8	1.74	0.69
2:I:22:DG:H1	3:J:1:DC:H42	1.41	0.68
1:B:106:GLU:OE2	1:H:59:ARG:NH2	2.25	0.68
1:H:157:ARG:HB2	3:L:4:DA:H5'	1.76	0.68
2:I:1:DT:C5'	2:K:22:DG:O3'	2.42	0.67
1:H:63:GLU:O	1:H:66:PRO:HD2	1.94	0.66
2:E:22:DG:H1	3:F:1:DC:H42	1.41	0.66
1:B:218:TRP:O	1:B:222:MET:HG2	1.99	0.63
3:L:13:DC:H2''	3:L:14:DA:N7	2.15	0.62
1:G:218:TRP:O	1:G:222:MET:HG2	2.00	0.62
2:K:2:DT:N1	2:K:2:DT:O4'	2.33	0.62
1:A:157:ARG:HB2	3:D:4:DA:H5'	1.80	0.61
2:K:6:DG:C8	2:K:7:DT:H73	2.34	0.61
3:D:22:DA:C2'	3:F:1:DC:H5''	2.29	0.61
2:C:6:DG:C2'	2:C:7:DT:H5'	2.31	0.61
2:K:6:DG:C2'	2:K:7:DT:H5'	2.31	0.61
2:E:6:DG:C2'	2:E:7:DT:H5'	2.31	0.61
2:K:11:DG:H2''	2:K:12:DG:C8	2.35	0.61
1:A:218:TRP:O	1:A:222:MET:HG2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:18:DG:H1	3:F:5:DC:H42	1.50	0.60
1:H:218:TRP:O	1:H:222:MET:HG2	2.02	0.60
2:C:11:DG:H2''	2:C:12:DG:C8	2.37	0.59
2:I:6:DG:C2'	2:I:7:DT:H5'	2.32	0.59
2:I:11:DG:H2''	2:I:12:DG:C8	2.37	0.59
2:C:22:DG:C2'	2:E:1:DT:C5'	2.77	0.59
2:E:11:DG:H2''	2:E:12:DG:C8	2.37	0.59
3:D:13:DC:H2''	3:D:14:DA:N7	2.17	0.59
2:E:6:DG:C8	2:E:7:DT:H73	2.38	0.59
3:F:13:DC:H2''	3:F:14:DA:N7	2.18	0.58
1:G:68:PHE:HB3	1:G:80:LEU:HD22	1.86	0.58
1:B:70:ALA:HB1	1:H:45:VAL:HA	1.86	0.58
1:A:118:LYS:HA	1:A:121:LEU:HD12	1.85	0.57
2:C:1:DT:C5'	2:E:22:DG:C2'	2.76	0.57
1:H:68:PHE:HB3	1:H:80:LEU:HD22	1.87	0.57
1:H:227:ARG:NH2	1:H:229:ASP:OD2	2.35	0.57
2:K:6:DG:N7	2:K:7:DT:H73	2.19	0.57
1:A:165:TYR:CE2	1:A:185:VAL:HG12	2.39	0.57
2:K:20:DT:H2''	2:K:21:DG:H8	1.69	0.57
1:B:227:ARG:NH2	1:B:229:ASP:OD2	2.38	0.57
1:H:165:TYR:CE2	1:H:185:VAL:HG12	2.40	0.57
3:J:13:DC:H2''	3:J:14:DA:N7	2.20	0.56
2:K:22:DG:O5'	2:K:22:DG:C8	2.53	0.56
2:C:6:DG:C8	2:C:7:DT:H73	2.40	0.56
2:E:6:DG:N7	2:E:7:DT:H73	2.19	0.56
1:A:63:GLU:O	1:A:66:PRO:HD2	2.05	0.56
1:G:227:ARG:NH2	1:G:229:ASP:OD2	2.35	0.56
1:B:165:TYR:CE2	1:B:185:VAL:HG12	2.40	0.56
3:D:22:DA:C3'	3:F:1:DC:H5''	2.36	0.56
3:F:6:DC:H2'	3:F:7:DA:C8	2.41	0.56
1:G:165:TYR:CE2	1:G:185:VAL:HG12	2.42	0.55
2:I:22:DG:C8	2:I:22:DG:O5'	2.55	0.55
2:I:21:DG:H2'	2:I:22:DG:N7	2.22	0.55
1:B:88:TRP:CZ2	1:B:96:LYS:HG2	2.42	0.54
2:I:6:DG:C8	2:I:7:DT:H73	2.42	0.54
1:G:88:TRP:CZ2	1:G:96:LYS:HG2	2.42	0.54
1:A:154:LYS:HB2	1:A:218:TRP:CZ2	2.42	0.54
1:H:88:TRP:CZ2	1:H:96:LYS:HG2	2.42	0.54
1:B:65:LEU:HB3	1:B:66:PRO:HD3	1.90	0.54
1:B:106:GLU:CD	1:H:59:ARG:NH2	2.62	0.53
3:J:1:DC:C5'	3:L:22:DA:H2''	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:63:GLU:C	1:H:66:PRO:HD2	2.29	0.53
1:A:65:LEU:HB3	1:A:66:PRO:HD3	1.90	0.53
3:J:6:DC:H2'	3:J:7:DA:C8	2.44	0.53
1:A:227:ARG:NH2	1:A:229:ASP:OD2	2.38	0.53
2:C:20:DT:H2''	2:C:21:DG:H8	1.73	0.53
1:B:118:LYS:HA	1:B:121:LEU:HD12	1.90	0.52
1:B:68:PHE:HB3	1:B:80:LEU:HD22	1.92	0.52
2:E:20:DT:H2''	2:E:21:DG:H8	1.74	0.52
1:H:65:LEU:HB3	1:H:66:PRO:HD3	1.92	0.52
3:J:2:DC:H2''	3:J:3:DA:C8	2.45	0.51
3:D:6:DC:H2'	3:D:7:DA:C8	2.45	0.51
3:J:2:DC:H2''	3:J:3:DA:H8	1.76	0.51
2:C:6:DG:N7	2:C:7:DT:H73	2.25	0.51
1:H:218:TRP:NE1	1:H:222:MET:HE3	2.25	0.51
2:K:20:DT:H2''	2:K:21:DG:C8	2.46	0.51
1:A:63:GLU:C	1:A:66:PRO:HD2	2.32	0.50
2:I:20:DT:H2''	2:I:21:DG:H8	1.77	0.50
1:B:225:VAL:O	1:B:225:VAL:HG12	2.11	0.50
1:H:118:LYS:HA	1:H:121:LEU:HD12	1.93	0.50
1:G:118:LYS:HA	1:G:121:LEU:HD12	1.94	0.49
1:G:158:PRO:HD3	1:G:211:TYR:CG	2.47	0.49
2:C:22:DG:O5'	2:C:22:DG:C8	2.63	0.49
1:B:79:GLU:CD	1:B:82:ARG:HH11	2.15	0.49
1:G:154:LYS:HB2	1:G:218:TRP:CZ2	2.47	0.49
1:A:165:TYR:HE2	1:A:185:VAL:HG12	1.78	0.49
3:D:2:DC:H2''	3:D:3:DA:C8	2.48	0.49
2:I:6:DG:N7	2:I:7:DT:H73	2.27	0.49
1:A:152:LEU:HB3	1:A:222:MET:HE3	1.95	0.49
1:H:165:TYR:HE2	1:H:185:VAL:HG12	1.78	0.49
3:J:22:DA:H2''	3:L:1:DC:C5'	2.43	0.49
3:L:6:DC:H2'	3:L:7:DA:C8	2.48	0.48
1:B:165:TYR:HE2	1:B:185:VAL:HG12	1.77	0.48
2:K:18:DG:H1	3:L:5:DC:H42	1.60	0.48
1:B:167:ALA:HA	3:F:6:DC:H5'	1.95	0.48
2:C:6:DG:H2''	2:C:7:DT:H5'	1.95	0.48
2:E:22:DG:C8	2:E:22:DG:O5'	2.57	0.48
1:B:158:PRO:HD3	1:B:211:TYR:CG	2.48	0.48
2:K:21:DG:H2'	2:K:22:DG:N7	2.29	0.48
2:I:6:DG:H2''	2:I:7:DT:H5'	1.95	0.48
1:H:158:PRO:HD3	1:H:211:TYR:CG	2.49	0.48
1:G:165:TYR:HE2	1:G:185:VAL:HG12	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:232:ARG:O	1:H:233:ARG:HB2	2.14	0.47
1:A:88:TRP:CZ2	1:A:96:LYS:HG2	2.48	0.47
2:I:11:DG:H2''	2:I:12:DG:H8	1.79	0.47
2:C:21:DG:H2'	2:C:22:DG:N7	2.30	0.47
3:F:16:DA:H62	3:F:17:DC:N4	2.13	0.47
2:K:11:DG:H2''	2:K:12:DG:H8	1.76	0.47
1:H:49:CYS:SG	1:H:50:PRO:HD2	2.55	0.46
2:C:18:DG:H1	3:D:5:DC:H42	1.63	0.46
3:D:2:DC:H2''	3:D:3:DA:H8	1.81	0.46
2:E:20:DT:H2''	2:E:21:DG:C8	2.51	0.46
2:K:6:DG:C8	2:K:7:DT:C7	2.98	0.46
2:E:6:DG:H2''	2:E:7:DT:H5'	1.97	0.46
2:C:20:DT:H2''	2:C:21:DG:C8	2.50	0.46
1:H:225:VAL:O	1:H:225:VAL:HG12	2.16	0.46
1:B:149:LEU:HD21	1:B:230:LEU:HD22	1.98	0.46
1:A:114:ILE:O	1:A:117:PHE:HB3	2.16	0.46
1:G:79:GLU:CD	1:G:82:ARG:HH11	2.14	0.46
1:A:124:SER:HB3	1:B:124:SER:HB3	1.99	0.46
1:A:64:GLN:O	1:A:65:LEU:C	2.54	0.45
1:B:207:ASP:HA	1:B:210:ARG:HE	1.81	0.45
3:D:8:DA:H2''	3:D:9:DG:N7	2.31	0.45
1:B:152:LEU:HB3	1:B:222:MET:HE3	1.97	0.45
1:B:59:ARG:NH2	1:H:106:GLU:OE2	2.49	0.45
2:E:21:DG:H2'	2:E:22:DG:N7	2.32	0.45
1:A:68:PHE:HB3	1:A:80:LEU:HD22	1.98	0.45
3:L:2:DC:H2''	3:L:3:DA:C8	2.51	0.45
1:B:113:GLU:OE2	1:H:63:GLU:HG3	2.16	0.45
3:J:8:DA:H2''	3:J:9:DG:N7	2.32	0.45
3:F:8:DA:H2''	3:F:9:DG:N7	2.32	0.45
1:G:63:GLU:C	1:G:66:PRO:HD2	2.36	0.45
2:C:1:DT:O4'	2:E:22:DG:H2''	2.17	0.45
1:B:65:LEU:HA	1:B:65:LEU:HD12	1.86	0.44
2:C:1:DT:C5'	2:E:22:DG:C3'	2.95	0.44
3:J:7:DA:H5''	3:J:7:DA:H8	1.82	0.44
1:A:158:PRO:HD3	1:A:211:TYR:CG	2.52	0.44
1:G:218:TRP:NE1	1:G:222:MET:HE3	2.32	0.44
1:G:207:ASP:HA	1:G:210:ARG:HE	1.83	0.44
2:E:6:DG:C8	2:E:7:DT:C7	3.00	0.44
3:L:8:DA:H2''	3:L:9:DG:N7	2.31	0.44
1:H:84:ILE:O	1:H:87:ARG:HB2	2.17	0.44
2:E:11:DG:H2''	2:E:12:DG:H8	1.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:22:DG:C3'	2:E:1:DT:C5'	2.96	0.44
1:A:177:SER:HB2	1:A:178:PRO:CD	2.47	0.44
2:I:1:DT:C5'	2:K:22:DG:C3'	2.96	0.44
2:I:20:DT:H2''	2:I:21:DG:C8	2.53	0.44
1:G:65:LEU:HA	1:G:65:LEU:HD12	1.74	0.44
2:C:22:DG:H2''	2:E:1:DT:O4'	2.18	0.43
3:L:7:DA:H5''	3:L:7:DA:H8	1.82	0.43
1:G:228:LYS:HG3	1:H:226:GLY:HA3	2.00	0.43
3:F:2:DC:H2''	3:F:3:DA:C8	2.53	0.43
3:F:16:DA:N7	3:F:17:DC:C4	2.87	0.43
3:F:2:DC:H2''	3:F:3:DA:H8	1.84	0.43
2:K:6:DG:H2'	2:K:7:DT:H5'	1.99	0.43
1:B:63:GLU:C	1:B:66:PRO:HD2	2.39	0.43
1:A:225:VAL:HG12	1:A:225:VAL:O	2.18	0.43
1:B:154:LYS:HB2	1:B:218:TRP:CZ2	2.54	0.43
2:K:6:DG:H2''	2:K:7:DT:H5'	2.00	0.43
1:H:149:LEU:HD21	1:H:230:LEU:HD22	2.01	0.43
2:C:11:DG:H2''	2:C:12:DG:H8	1.80	0.42
1:G:225:VAL:HG12	1:G:225:VAL:O	2.19	0.42
1:B:99:TYR:O	1:B:102:ALA:HB3	2.18	0.42
3:L:2:DC:H2''	3:L:3:DA:H8	1.85	0.42
3:L:8:DA:H2''	3:L:9:DG:C8	2.55	0.42
1:G:63:GLU:O	1:G:66:PRO:HD2	2.20	0.42
1:H:207:ASP:HA	1:H:210:ARG:HE	1.85	0.42
1:B:77:THR:O	1:B:81:ILE:HG12	2.19	0.42
1:H:114:ILE:HG23	1:H:115:SER:N	2.34	0.42
3:J:8:DA:H2''	3:J:9:DG:C8	2.55	0.41
2:C:6:DG:C8	2:C:7:DT:C7	3.03	0.41
2:E:16:DT:C4	2:E:17:DG:O6	2.73	0.41
1:G:49:CYS:SG	1:G:50:PRO:HD2	2.60	0.41
1:A:157:ARG:HH11	1:A:157:ARG:HG3	1.85	0.41
2:I:22:DG:C3'	2:K:1:DT:C5'	2.98	0.41
1:B:114:ILE:O	1:B:117:PHE:HB3	2.19	0.41
1:H:65:LEU:HA	1:H:65:LEU:HD12	1.87	0.41
1:H:114:ILE:O	1:H:117:PHE:HB3	2.21	0.41
2:C:16:DT:C4	2:C:17:DG:O6	2.74	0.41
1:H:154:LYS:HB2	1:H:218:TRP:CZ2	2.55	0.41
1:B:116:ARG:HH21	1:H:64:GLN:NE2	2.18	0.41
1:G:59:ARG:O	1:G:60:PHE:C	2.59	0.41
1:B:52:LYS:NZ	3:F:22:DA:OP1	2.49	0.41
3:D:16:DA:H62	3:D:17:DC:N4	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:154:LYS:HD2	1:H:218:TRP:CH2	2.56	0.41
3:F:8:DA:H2''	3:F:9:DG:C8	2.55	0.41
1:G:177:SER:HB2	1:G:178:PRO:CD	2.51	0.41
3:D:8:DA:H2''	3:D:9:DG:C8	2.55	0.40
2:I:16:DT:C4	2:I:17:DG:O6	2.73	0.40
2:C:6:DG:H2'	2:C:7:DT:H5'	2.03	0.40
1:H:177:SER:HB2	1:H:178:PRO:CD	2.51	0.40
1:B:177:SER:HB2	1:B:178:PRO:CD	2.51	0.40
1:G:55:SER:HA	1:G:103:TYR:CD1	2.57	0.40
1:H:58:LEU:HD11	2:K:6:DG:C2	2.56	0.40
1:A:59:ARG:O	1:A:60:PHE:C	2.59	0.40
3:F:15:DT:H2''	3:F:16:DA:H5'	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	188/230 (82%)	171 (91%)	17 (9%)	0	100	100
1	B	188/230 (82%)	173 (92%)	15 (8%)	0	100	100
1	G	188/230 (82%)	173 (92%)	15 (8%)	0	100	100
1	H	188/230 (82%)	174 (93%)	14 (7%)	0	100	100
All	All	752/920 (82%)	691 (92%)	61 (8%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar

resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	176/208 (85%)	175 (99%)	1 (1%)	90	97
1	B	176/208 (85%)	175 (99%)	1 (1%)	90	97
1	G	176/208 (85%)	175 (99%)	1 (1%)	90	97
1	H	176/208 (85%)	175 (99%)	1 (1%)	90	97
All	All	704/832 (85%)	700 (99%)	4 (1%)	90	97

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	150	THR
1	B	150	THR
1	G	150	THR
1	H	150	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	HIS
1	B	137	HIS
1	G	137	HIS
1	H	137	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected

value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	BRU	C	19	3,2	13,21,22	2.47	1 (7%)	16,30,33	2.33	2 (12%)
2	BRU	E	19	3,2	13,21,22	2.40	1 (7%)	16,30,33	2.28	2 (12%)
2	BRU	I	19	3,2	13,21,22	2.45	1 (7%)	16,30,33	2.35	2 (12%)
2	BRU	K	19	3,2	13,21,22	2.41	1 (7%)	16,30,33	2.25	2 (12%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BRU	C	19	3,2	-	0/3/21/22	0/2/2/2
2	BRU	E	19	3,2	-	0/3/21/22	0/2/2/2
2	BRU	I	19	3,2	-	0/3/21/22	0/2/2/2
2	BRU	K	19	3,2	-	0/3/21/22	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	K	19	BRU	C4-C5	8.53	1.49	1.38
2	E	19	BRU	C4-C5	8.54	1.49	1.38
2	I	19	BRU	C4-C5	8.68	1.49	1.38
2	C	19	BRU	C4-C5	8.81	1.49	1.38

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	19	BRU	C5-C4-N3	-5.39	118.24	124.00
2	K	19	BRU	C5-C4-N3	-5.31	118.33	124.00
2	C	19	BRU	C5-C4-N3	-5.24	118.40	124.00
2	E	19	BRU	C5-C4-N3	-5.07	118.58	124.00
2	K	19	BRU	C4-N3-C2	6.70	121.03	115.25
2	I	19	BRU	C4-N3-C2	6.91	121.22	115.25
2	C	19	BRU	C4-N3-C2	7.05	121.34	115.25
2	E	19	BRU	C4-N3-C2	7.12	121.40	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

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	190/230 (82%)	0.18	17 (8%) 12 7	47, 92, 187, 227	0
1	B	190/230 (82%)	0.13	13 (6%) 20 14	43, 79, 179, 206	0
1	G	190/230 (82%)	-0.21	1 (0%) 91 90	42, 70, 140, 153	0
1	H	190/230 (82%)	-0.17	1 (0%) 91 90	38, 78, 140, 162	0
2	C	21/22 (95%)	-0.30	0 100 100	65, 102, 177, 185	0
2	E	21/22 (95%)	-0.31	0 100 100	69, 111, 174, 205	0
2	I	21/22 (95%)	-0.26	0 100 100	54, 81, 155, 166	0
2	K	21/22 (95%)	-0.28	0 100 100	57, 81, 154, 181	0
3	D	22/22 (100%)	-0.08	0 100 100	62, 105, 201, 243	0
3	F	22/22 (100%)	-0.12	1 (4%) 37 31	69, 105, 218, 247	0
3	J	22/22 (100%)	-0.15	0 100 100	53, 73, 176, 227	0
3	L	22/22 (100%)	-0.26	0 100 100	47, 77, 181, 240	0
All	All	932/1096 (85%)	-0.05	33 (3%) 48 40	38, 81, 178, 247	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	199	LEU	8.9
1	B	199	LEU	7.3
1	A	171	GLN	6.4
1	B	195	SER	5.3
1	A	161	ALA	5.2
1	B	176	ASP	5.1
1	A	164	VAL	5.0
1	A	189	TRP	4.3
1	B	171	GLN	4.2
3	F	10	DC	3.8
1	B	174	LYS	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	172	GLU	3.7
1	A	176	ASP	3.3
1	A	200	TYR	3.1
1	B	200	TYR	3.0
1	A	191	ASN	3.0
1	A	172	GLU	2.9
1	H	165	TYR	2.9
1	B	196	GLU	2.8
1	A	165	TYR	2.8
1	A	182	LEU	2.6
1	B	192	LEU	2.6
1	A	198	GLU	2.5
1	B	191	ASN	2.5
1	A	188	ASN	2.5
1	A	205	LYS	2.4
1	A	184	THR	2.3
1	G	169	ARG	2.2
1	A	203	HIS	2.2
1	A	195	SER	2.1
1	B	184	THR	2.1
1	B	165	TYR	2.0
1	B	175	GLY	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	BRU	E	19	20/21	0.91	0.09	-	135,141,179,180	0
2	BRU	I	19	20/21	0.95	0.11	-	95,109,126,147	0
2	BRU	K	19	20/21	0.96	0.08	-	104,115,120,158	0
2	BRU	C	19	20/21	0.88	0.10	-	120,142,155,160	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.