



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:49 AM GMT

PDB ID : 3SWP  
Title : ANAC019 NAC domain in complex with DNA  
Authors : Welner, D.; Lo Leggio, L.  
Deposited on : 2011-07-14  
Resolution : 4.11 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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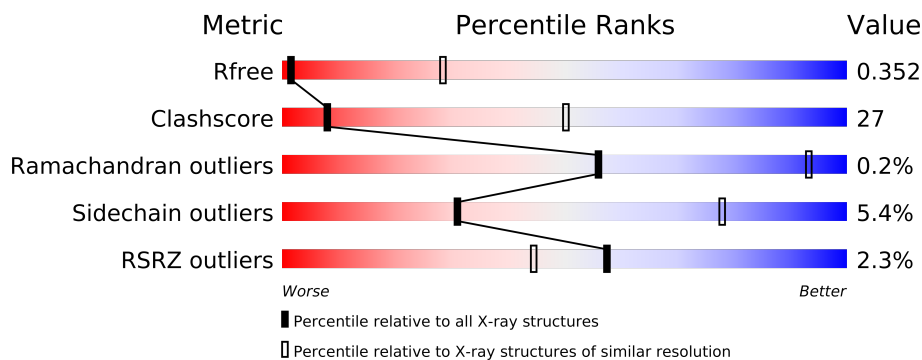
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 4.11 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))
$R_{free}$	66092	1003 (4.76-3.48)
Clashscore	79885	1248 (4.70-3.50)
Ramachandran outliers	78287	1183 (4.70-3.50)
Sidechain outliers	78261	1168 (4.70-3.50)
RSRZ outliers	66119	1003 (4.76-3.48)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	174	
1	B	174	
1	C	174	
1	D	174	
2	E	26	
3	F	26	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5667 atoms, of which 0 are hydrogen and 0 are deuterium.

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 NAC domain-containing protein 19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	0	0	0
			1167	767	193	203	4			
1	B	140	Total	C	N	O	S	0	0	0
			1156	761	189	202	4			
1	C	135	Total	C	N	O	S	0	0	0
			1122	738	186	194	4			
1	D	140	Total	C	N	O	S	0	0	0
			1156	761	189	202	4			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	EXPRESSION TAG	UNP Q9C932
A	-4	HIS	-	EXPRESSION TAG	UNP Q9C932
A	-3	HIS	-	EXPRESSION TAG	UNP Q9C932
A	-2	HIS	-	EXPRESSION TAG	UNP Q9C932
A	-1	HIS	-	EXPRESSION TAG	UNP Q9C932
A	0	HIS	-	EXPRESSION TAG	UNP Q9C932
B	-5	HIS	-	EXPRESSION TAG	UNP Q9C932
B	-4	HIS	-	EXPRESSION TAG	UNP Q9C932
B	-3	HIS	-	EXPRESSION TAG	UNP Q9C932
B	-2	HIS	-	EXPRESSION TAG	UNP Q9C932
B	-1	HIS	-	EXPRESSION TAG	UNP Q9C932
B	0	HIS	-	EXPRESSION TAG	UNP Q9C932
C	-5	HIS	-	EXPRESSION TAG	UNP Q9C932
C	-4	HIS	-	EXPRESSION TAG	UNP Q9C932
C	-3	HIS	-	EXPRESSION TAG	UNP Q9C932
C	-2	HIS	-	EXPRESSION TAG	UNP Q9C932
C	-1	HIS	-	EXPRESSION TAG	UNP Q9C932
C	0	HIS	-	EXPRESSION TAG	UNP Q9C932
D	-5	HIS	-	EXPRESSION TAG	UNP Q9C932
D	-4	HIS	-	EXPRESSION TAG	UNP Q9C932
D	-3	HIS	-	EXPRESSION TAG	UNP Q9C932

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	HIS	-	EXPRESSION TAG	UNP Q9C932
D	-1	HIS	-	EXPRESSION TAG	UNP Q9C932
D	0	HIS	-	EXPRESSION TAG	UNP Q9C932

- Molecule 2 is a DNA chain called oligonucleotide forward.


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	26	Total	C	N	O	P	0	0	0
			536	254	100	156	26			

- Molecule 3 is a DNA chain called oligonucleotide reverse.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	26	Total	C	N	O	P	0	0	0
			530	252	96	156	26			




● Molecule 2: oligonucleotide forward

Chain E: 

G1	T2	C3	T4	T5	G6	C7	G8	T9	G10	T11	T12	G13	G14	A15	A16	C17	A18	C19	G20	C21	A22	A23	C24	A25	G26
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● Molecule 3: oligonucleotide reverse

Chain F: 

C1	C2	T3	G4	T5	T6	G7	C8	G9	T10	G11	T12	T13	C14	C15	A16	A17	C18	A19	C20	G21	C22	A23	A24	G25	A26
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	69.14Å 105.47Å 175.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.97 – 4.11 29.97 – 4.11	Depositor EDS
% Data completeness (in resolution range)	99.0 (29.97-4.11) 99.2 (29.97-4.11)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.05 (at 4.11Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.289)	Depositor
R, $R_{free}$	0.260 , 0.349 0.257 , 0.352	Depositor DCC
$R_{free}$ test set	1035 reflections (10.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	174.1	Xtriage
Anisotropy	0.465	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.22 , 140.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 10348 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	5667	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	263.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.41% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 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	0.94	1/1202 (0.1%)	1.40	13/1623 (0.8%)
1	B	0.62	1/1191 (0.1%)	1.09	8/1609 (0.5%)
1	C	0.87	1/1156 (0.1%)	1.38	14/1560 (0.9%)
1	D	0.58	1/1191 (0.1%)	1.06	8/1609 (0.5%)
2	E	0.89	2/601 (0.3%)	1.84	25/926 (2.7%)
3	F	0.82	0/593	2.08	35/912 (3.8%)
All	All	0.79	6/5934 (0.1%)	1.43	103/8239 (1.3%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	10	THR	CA-C	6.29	1.69	1.52
1	B	10	THR	CA-C	6.26	1.69	1.52
2	E	6	DG	C3'-O3'	-5.58	1.36	1.44
1	C	157	CYS	CB-SG	-5.41	1.73	1.81
1	A	132	TRP	CB-CG	-5.21	1.40	1.50
2	E	7	DC	C1'-N1	5.20	1.56	1.49

All (103) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	19	DA	C1'-O4'-C4'	-12.87	97.23	110.10
3	F	19	DA	O4'-C1'-N9	10.88	115.62	108.00
1	C	45	LEU	CA-CB-CG	-9.44	93.59	115.30
1	A	45	LEU	CA-CB-CG	-9.39	93.71	115.30
3	F	5	DT	O4'-C1'-N1	9.30	114.51	108.00
2	E	12	DT	O4'-C1'-N1	-8.80	101.84	108.00
2	E	9	DT	O4'-C1'-N1	8.68	114.08	108.00
3	F	13	DT	O4'-C1'-C2'	-8.54	99.07	105.90
3	F	26	DA	C4'-C3'-C2'	-8.49	95.46	103.10
2	E	10	DG	O4'-C4'-C3'	-8.29	101.02	106.00
3	F	9	DG	O4'-C1'-C2'	-8.29	99.26	105.90
1	A	67	GLU	N-CA-CB	-8.22	95.80	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	11	DG	O4'-C1'-C2'	-8.16	99.37	105.90
3	F	2	DC	O4'-C1'-N1	8.15	113.71	108.00
1	C	67	GLU	N-CA-CB	-7.95	96.30	110.60
2	E	21	DC	O4'-C4'-C3'	-7.89	101.27	106.00
3	F	26	DA	O4'-C1'-C2'	-7.86	99.61	105.90
3	F	26	DA	O4'-C1'-N9	7.80	113.46	108.00
3	F	19	DA	C3'-C2'-C1'	-7.47	93.54	102.50
3	F	15	DC	O4'-C1'-N1	7.37	113.16	108.00
2	E	23	DA	C3'-C2'-C1'	-7.32	93.71	102.50
1	D	11	GLN	N-CA-CB	7.20	123.55	110.60
3	F	3	DT	O4'-C4'-C3'	-7.18	101.63	104.50
1	C	32	LEU	CA-CB-CG	7.13	131.70	115.30
2	E	15	DA	O4'-C1'-N9	7.09	112.96	108.00
1	B	11	GLN	N-CA-CB	7.06	123.31	110.60
1	A	32	LEU	CA-CB-CG	7.05	131.52	115.30
1	C	49	ILE	CG1-CB-CG2	-7.05	95.89	111.40
2	E	26	DG	C4'-C3'-C2'	-7.01	96.79	103.10
3	F	13	DT	O4'-C1'-N1	6.96	112.87	108.00
1	C	117	LEU	CB-CG-CD1	-6.94	99.19	111.00
2	E	12	DT	C4-C5-C7	6.84	123.10	119.00
3	F	10	DT	O4'-C1'-C2'	-6.81	100.45	105.90
2	E	17	DC	O4'-C1'-N1	6.77	112.74	108.00
1	A	117	LEU	CB-CG-CD1	-6.76	99.50	111.00
3	F	6	DT	O4'-C1'-N1	6.75	112.72	108.00
3	F	12	DT	C4'-C3'-C2'	-6.66	97.11	103.10
3	F	14	DC	O4'-C1'-N1	6.61	112.62	108.00
2	E	5	DT	N3-C4-O4	6.50	123.80	119.90
2	E	20	DG	C3'-C2'-C1'	-6.48	94.73	102.50
1	B	21	TYR	CA-CB-CG	6.38	125.53	113.40
1	D	21	TYR	CA-CB-CG	6.31	125.40	113.40
2	E	9	DT	C3'-C2'-C1'	-6.31	94.93	102.50
1	C	156	LEU	CB-CG-CD1	-6.28	100.33	111.00
1	A	28	MET	CA-CB-CG	6.27	123.95	113.30
1	A	49	ILE	CG1-CB-CG2	-6.22	97.71	111.40
3	F	11	DG	C3'-C2'-C1'	-6.21	95.04	102.50
2	E	11	DT	O4'-C1'-N1	6.13	112.29	108.00
3	F	19	DA	P-O5'-C5'	-6.12	111.10	120.90
3	F	5	DT	O4'-C1'-C2'	-6.09	101.03	105.90
1	A	59	LEU	N-CA-C	6.08	127.42	111.00
1	A	162	LYS	N-CA-C	6.08	127.41	111.00
1	D	10	THR	CB-CA-C	-6.04	95.28	111.60
1	B	10	THR	CB-CA-C	-6.01	95.36	111.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	13	DT	C1'-O4'-C4'	-5.92	104.19	110.10
3	F	10	DT	C1'-O4'-C4'	-5.91	104.19	110.10
3	F	26	DA	C3'-C2'-C1'	-5.89	95.43	102.50
1	C	162	LYS	N-CA-C	5.86	126.82	111.00
1	C	28	MET	CA-CB-CG	5.82	123.19	113.30
3	F	10	DT	O4'-C1'-N1	5.76	112.03	108.00
1	C	59	LEU	N-CA-C	5.70	126.40	111.00
3	F	10	DT	C3'-C2'-C1'	-5.68	95.69	102.50
1	B	10	THR	CA-C-O	5.65	131.97	120.10
2	E	5	DT	C5-C4-O4	-5.62	120.97	124.90
3	F	19	DA	C8-N9-C4	5.62	108.05	105.80
1	D	10	THR	CA-C-O	5.55	131.75	120.10
1	C	33	CYS	CA-CB-SG	-5.51	104.08	114.00
3	F	7	DG	C3'-C2'-C1'	-5.48	95.92	102.50
1	A	22	PRO	N-CA-C	-5.48	97.86	112.10
1	A	156	LEU	CB-CG-CD1	-5.47	101.69	111.00
1	A	100	THR	N-CA-C	-5.44	96.32	111.00
1	B	10	THR	CA-C-N	-5.43	105.26	117.20
3	F	6	DT	C1'-O4'-C4'	-5.41	104.69	110.10
1	C	100	THR	N-CA-C	-5.41	96.41	111.00
2	E	21	DC	C3'-C2'-C1'	-5.39	96.03	102.50
1	C	22	PRO	N-CA-C	-5.38	98.11	112.10
1	A	162	LYS	CB-CA-C	-5.38	99.65	110.40
1	D	10	THR	CA-C-N	-5.34	105.46	117.20
3	F	5	DT	N3-C4-O4	5.33	123.09	119.90
2	E	13	DG	C3'-C2'-C1'	-5.29	96.15	102.50
3	F	10	DT	N3-C4-O4	5.29	123.08	119.90
3	F	5	DT	C1'-O4'-C4'	-5.29	104.81	110.10
2	E	16	DA	C4'-C3'-C2'	-5.27	98.36	103.10
2	E	12	DT	C6-C5-C7	-5.26	119.74	122.90
3	F	15	DC	C1'-O4'-C4'	-5.17	104.93	110.10
1	B	21	TYR	N-CA-CB	5.16	119.89	110.60
3	F	12	DT	N3-C4-O4	5.15	122.99	119.90
2	E	2	DT	O4'-C1'-N1	5.15	111.61	108.00
1	C	162	LYS	CB-CA-C	-5.15	100.10	110.40
1	B	64	LEU	CA-CB-CG	5.13	127.11	115.30
2	E	2	DT	N3-C4-O4	5.12	122.97	119.90
3	F	14	DC	C1'-O4'-C4'	-5.12	104.98	110.10
2	E	1	DG	O4'-C4'-C3'	-5.11	102.46	104.50
1	D	64	LEU	CA-CB-CG	5.10	127.04	115.30
2	E	20	DG	O4'-C1'-C2'	-5.10	101.82	105.90
1	B	11	GLN	N-CA-C	-5.08	97.29	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	ILE	CB-CA-C	-5.05	101.49	111.60
2	E	23	DA	C1'-O4'-C4'	-5.05	105.05	110.10
1	C	76	ARG	N-CA-C	5.04	124.61	111.00
1	D	10	THR	N-CA-C	5.04	124.61	111.00
1	D	21	TYR	N-CA-CB	5.04	119.67	110.60
2	E	2	DT	C5-C4-O4	-5.02	121.38	124.90
2	E	9	DT	O4'-C1'-C2'	-5.02	101.88	105.90

There are no chirality outliers.

There are no planarity outliers.

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1167	0	0	25	0
1	B	1156	0	0	12	0
1	C	1122	0	0	21	0
1	D	1156	0	0	16	0
2	E	536	0	0	51	0
3	F	530	0	0	58	0
All	All	5667	0	0	143	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 27.

All (143) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:21:TYR:CD2	1:D:21:TYR:CD2	2.43	1.06
1:A:129:LYS:NZ	3:F:20:DC:P	2.31	1.03
1:A:129:LYS:NZ	3:F:20:DC:O5'	1.95	0.97
1:D:87:ASN:ND2	3:F:12:DT:OP2	2.04	0.91
1:A:129:LYS:NZ	3:F:20:DC:OP2	2.04	0.90
3:F:17:DA:N7	3:F:18:DC:C4	2.40	0.89
2:E:5:DT:C2'	2:E:6:DG:C8	2.59	0.85

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:96:LYS:NZ	3:F:8:DC:N4	2.27	0.82
1:C:135:HIS:ND1	1:C:158:ARG:NH1	2.30	0.79
3:F:17:DA:C8	3:F:18:DC:C4	2.72	0.77
2:E:5:DT:O4	3:F:23:DA:N1	2.18	0.77
2:E:9:DT:O2	2:E:10:DG:C5	2.38	0.76
3:F:11:DG:N7	3:F:12:DT:O4	2.21	0.73
2:E:10:DG:C2	2:E:11:DT:C4	2.77	0.73
1:A:21:TYR:CD2	1:B:21:TYR:CD2	2.77	0.72
3:F:17:DA:N7	3:F:18:DC:N3	2.38	0.71
1:A:135:HIS:ND1	1:A:158:ARG:NH1	2.39	0.70
2:E:18:DA:N1	3:F:10:DT:O4	2.25	0.70
1:C:21:TYR:CG	1:D:21:TYR:CE2	2.79	0.70
3:F:11:DG:N7	3:F:12:DT:C4	2.61	0.68
1:D:86:PRO:O	1:D:87:ASN:C	2.32	0.68
1:C:19:ARG:CZ	1:D:26:GLU:OE2	2.41	0.68
2:E:14:DG:N2	3:F:14:DC:N3	2.43	0.67
3:F:25:DG:C1'	3:F:26:DA:C8	2.79	0.66
1:B:86:PRO:O	1:B:87:ASN:C	2.33	0.66
1:C:21:TYR:CB	1:D:21:TYR:CE2	2.81	0.63
2:E:13:DG:O6	3:F:15:DC:N4	2.32	0.62
1:A:36:ALA:C	1:A:38:GLY:N	2.52	0.62
1:B:129:LYS:CE	2:E:17:DC:OP2	2.48	0.62
2:E:10:DG:C2'	2:E:11:DT:OP2	2.47	0.62
1:C:19:ARG:NH2	1:D:26:GLU:OE2	2.33	0.62
1:A:135:HIS:O	1:A:158:ARG:N	2.34	0.60
3:F:20:DC:C4	3:F:21:DG:N1	2.69	0.60
1:C:21:TYR:CG	1:D:21:TYR:CD2	2.89	0.60
2:E:8:DG:N1	3:F:19:DA:C6	2.69	0.60
3:F:13:DT:C2'	3:F:14:DC:C6	2.84	0.60
1:A:12:LEU:O	1:A:13:SER:C	2.41	0.59
2:E:8:DG:C6	3:F:19:DA:N6	2.70	0.59
2:E:14:DG:C2'	2:E:15:DA:O4'	2.49	0.59
1:A:86:PRO:CB	2:E:4:DT:C4'	2.80	0.59
1:C:63:ALA:CB	1:C:70:TRP:CD1	2.85	0.59
1:C:36:ALA:C	1:C:38:GLY:N	2.55	0.58
2:E:9:DT:O2	2:E:10:DG:N7	2.36	0.58
2:E:17:DC:C5	2:E:18:DA:N6	2.71	0.58
1:A:28:MET:O	1:A:32:LEU:CB	2.51	0.57
2:E:5:DT:O4	3:F:23:DA:C6	2.56	0.57
1:A:86:PRO:O	2:E:4:DT:OP1	2.21	0.57
2:E:8:DG:C2	2:E:9:DT:N3	2.73	0.57
1:B:129:LYS:NZ	2:E:16:DA:C3'	2.68	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:129:LYS:CE	3:F:20:DC:OP2	2.52	0.56
2:E:10:DG:N2	2:E:11:DT:N3	2.54	0.56
3:F:17:DA:C8	3:F:18:DC:C5	2.94	0.56
1:C:28:MET:O	1:C:32:LEU:CB	2.53	0.56
3:F:11:DG:C8	3:F:12:DT:C4	2.94	0.56
1:C:135:HIS:O	1:C:158:ARG:N	2.39	0.55
1:A:63:ALA:CB	1:A:70:TRP:CD1	2.90	0.55
3:F:12:DT:C2'	3:F:13:DT:C6	2.90	0.55
3:F:25:DG:N3	3:F:26:DA:C8	2.75	0.55
3:F:25:DG:N3	3:F:26:DA:C5	2.75	0.55
1:D:162:LYS:NZ	2:E:11:DT:OP1	2.40	0.55
3:F:20:DC:N3	3:F:21:DG:C2	2.75	0.54
3:F:23:DA:C5	3:F:24:DA:C6	2.95	0.54
1:B:23:THR:OG1	1:B:26:GLU:N	2.40	0.54
3:F:20:DC:C4	3:F:21:DG:C2	2.96	0.54
1:A:21:TYR:CD2	1:B:21:TYR:CE2	2.96	0.54
3:F:24:DA:C5	3:F:25:DG:O6	2.61	0.53
1:A:137:TYR:N	1:A:156:LEU:O	2.42	0.53
2:E:8:DG:C6	2:E:9:DT:O4	2.61	0.53
2:E:26:DG:N1	3:F:2:DC:N4	2.56	0.53
1:D:23:THR:OG1	1:D:26:GLU:N	2.42	0.52
2:E:18:DA:C2	3:F:11:DG:N2	2.78	0.52
2:E:8:DG:C2	2:E:9:DT:C4	2.98	0.52
1:A:57:TRP:CH2	1:A:121:ILE:CD1	2.93	0.52
1:C:137:TYR:N	1:C:156:LEU:O	2.42	0.51
3:F:8:DC:OP2	3:F:8:DC:C6	2.64	0.51
1:D:86:PRO:CB	3:F:12:DT:C4'	2.89	0.51
2:E:10:DG:N3	2:E:11:DT:C4	2.79	0.51
3:F:25:DG:C2	3:F:26:DA:C5	2.99	0.50
3:F:17:DA:N7	3:F:18:DC:N4	2.60	0.50
1:D:22:PRO:CA	1:D:26:GLU:OE1	2.61	0.49
1:A:86:PRO:CB	2:E:4:DT:C3'	2.90	0.49
1:B:10:THR:CG2	1:B:10:THR:O	2.59	0.49
3:F:25:DG:N2	3:F:26:DA:C2	2.80	0.49
2:E:26:DG:N2	3:F:2:DC:N3	2.61	0.49
1:A:35:LYS:O	1:A:38:GLY:CA	2.61	0.48
1:C:140:ILE:CG2	1:C:141:GLU:N	2.76	0.48
1:D:98:THR:OG1	2:E:12:DT:C7	2.62	0.48
2:E:10:DG:C2	2:E:11:DT:N3	2.82	0.48
1:B:22:PRO:CA	1:B:26:GLU:OE1	2.62	0.47
3:F:11:DG:C8	3:F:12:DT:C7	2.98	0.47
2:E:26:DG:O6	3:F:1:DC:N4	2.47	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:E:14:DG:O6	3:F:13:DT:O4	2.33	0.47
2:E:4:DT:C2'	2:E:5:DT:C6	2.98	0.47
1:C:34:ARG:O	1:C:39:TYR:O	2.33	0.47
2:E:9:DT:C6	2:E:9:DT:OP2	2.68	0.47
3:F:11:DG:C8	3:F:12:DT:C5	3.02	0.47
2:E:25:DA:C5	2:E:26:DG:C6	3.03	0.47
1:A:27:LEU:CD1	1:A:137:TYR:CE1	2.99	0.47
3:F:9:DG:C8	3:F:10:DT:C7	2.98	0.46
3:F:20:DC:C5	3:F:21:DG:C6	3.03	0.46
3:F:10:DT:O2	3:F:11:DG:C4	2.68	0.46
2:E:14:DG:N2	3:F:14:DC:C2	2.82	0.46
3:F:21:DG:C4	3:F:22:DC:N4	2.83	0.46
1:A:34:ARG:O	1:A:39:TYR:O	2.34	0.46
2:E:9:DT:C2	2:E:10:DG:C5	3.03	0.45
2:E:23:DA:C2'	2:E:24:DC:OP2	2.64	0.45
2:E:9:DT:C2	2:E:10:DG:N7	2.84	0.45
2:E:17:DC:C5	2:E:18:DA:C6	3.04	0.45
1:C:21:TYR:CD2	1:D:21:TYR:CE2	2.95	0.45
3:F:25:DG:C4'	3:F:26:DA:OP1	2.65	0.45
1:C:27:LEU:CD1	1:C:137:TYR:CE1	2.99	0.45
3:F:5:DT:C2'	3:F:6:DT:OP2	2.65	0.45
3:F:25:DG:N3	3:F:26:DA:C4	2.85	0.45
2:E:15:DA:N1	3:F:13:DT:O4	2.51	0.44
1:A:15:PRO:CB	1:A:18:PHE:CD2	3.00	0.44
1:A:35:LYS:NZ	1:A:48:GLU:OE2	2.50	0.44
2:E:8:DG:C6	3:F:19:DA:C6	3.05	0.44
1:C:57:TRP:CH2	1:C:121:ILE:CD1	3.00	0.44
1:D:25:GLU:OE2	1:D:106:THR:OG1	2.36	0.44
2:E:25:DA:C6	2:E:26:DG:C6	3.06	0.43
1:A:59:LEU:O	1:A:62:LYS:N	2.51	0.43
2:E:11:DT:C2'	2:E:11:DT:OP2	2.67	0.43
2:E:8:DG:C4	2:E:9:DT:C4	3.06	0.43
1:B:34:ARG:CD	1:B:39:TYR:CE2	3.01	0.43
2:E:23:DA:N6	3:F:5:DT:C4	2.87	0.42
1:A:33:CYS:O	1:A:34:ARG:C	2.57	0.42
2:E:14:DG:N2	3:F:14:DC:O2	2.52	0.42
3:F:12:DT:C6	3:F:12:DT:C5'	3.02	0.42
2:E:19:DC:C2'	2:E:20:DG:C8	3.02	0.42
1:B:59:LEU:N	1:B:60:PRO:CD	2.83	0.42
2:E:18:DA:N3	2:E:19:DC:C2	2.88	0.42
1:C:59:LEU:O	1:C:62:LYS:N	2.53	0.41
1:C:32:LEU:O	1:C:36:ALA:N	2.53	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:140:ILE:CG2	1:A:141:GLU:N	2.79	0.41
1:B:25:GLU:OE2	1:B:106:THR:OG1	2.38	0.41
2:E:18:DA:C4	2:E:19:DC:C4	3.09	0.40
3:F:25:DG:N2	3:F:26:DA:C4	2.90	0.40
1:D:59:LEU:N	1:D:60:PRO:CD	2.85	0.40
1:C:40:ASP:OD1	1:C:41:PHE:N	2.54	0.40
3:F:21:DG:C5	3:F:22:DC:N4	2.90	0.40
3:F:23:DA:C6	3:F:24:DA:N1	2.90	0.40
2:E:25:DA:C5	2:E:26:DG:O6	2.73	0.40
1:C:51:LEU:CD1	1:C:157:CYS:SG	3.10	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	135/174 (78%)	131 (97%)	3 (2%)	1 (1%)	30	84
1	B	134/174 (77%)	130 (97%)	4 (3%)	0	100	100
1	C	129/174 (74%)	127 (98%)	2 (2%)	0	100	100
1	D	134/174 (77%)	130 (97%)	4 (3%)	0	100	100
All	All	532/696 (76%)	518 (97%)	13 (2%)	1 (0%)	56	93

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	13	SER

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain

conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	123/152 (81%)	116 (94%)	7 (6%)	29	76
1	B	122/152 (80%)	116 (95%)	6 (5%)	35	80
1	C	117/152 (77%)	110 (94%)	7 (6%)	27	75
1	D	122/152 (80%)	116 (95%)	6 (5%)	35	80
All	All	484/608 (80%)	458 (95%)	26 (5%)	31	78

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	SER
1	A	53	LYS
1	A	140	ILE
1	A	143	SER
1	A	152	ASP
1	A	162	LYS
1	A	163	GLN
1	B	11	GLN
1	B	34	ARG
1	B	61	ASN
1	B	143	SER
1	B	162	LYS
1	B	163	GLN
1	C	42	SER
1	C	53	LYS
1	C	140	ILE
1	C	143	SER
1	C	152	ASP
1	C	162	LYS
1	C	163	GLN
1	D	11	GLN
1	D	34	ARG
1	D	61	ASN
1	D	143	SER
1	D	162	LYS
1	D	163	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.



### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates ⓘ

There are no 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	141/174 (81%)	-0.02	1 (0%) 84 73	127, 194, 305, 403	0
1	B	140/174 (80%)	0.31	7 (5%) 28 24	165, 269, 382, 579	0
1	C	135/174 (77%)	0.03	3 (2%) 59 46	137, 209, 331, 529	0
1	D	140/174 (80%)	0.17	3 (2%) 60 47	163, 274, 354, 560	0
2	E	26/26 (100%)	-0.24	0 100 100	249, 323, 393, 449	0
3	F	26/26 (100%)	-0.22	0 100 100	230, 321, 445, 481	0
All	All	608/748 (81%)	0.09	14 (2%) 57 44	127, 243, 384, 579	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	143	SER	9.1
1	B	152	ASP	4.9
1	B	142	PRO	4.4
1	C	78	ARG	3.1
1	D	142	PRO	2.8
1	A	110	ARG	2.8
1	C	14	LEU	2.3
1	B	131	ASN	2.3
1	D	157	CYS	2.3
1	B	48	GLU	2.3
1	B	77	ASP	2.3
1	C	110	ARG	2.3
1	B	141	GLU	2.3
1	D	143	SER	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

There are no ligands in this entry.

### 6.5 Other polymers ⓘ

There are no such residues in this entry.