



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 01:34 pm BST

PDB ID : 3A01  
Title : Crystal structure of Aristaless and Clawless homeodomains bound to DNA  
Authors : Miyazono, K.; Nagata, K.; Saigo, K.; Kojima, T.; Tanokura, M.  
Deposited on : 2009-02-28  
Resolution : 2.70 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](mailto: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.

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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
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

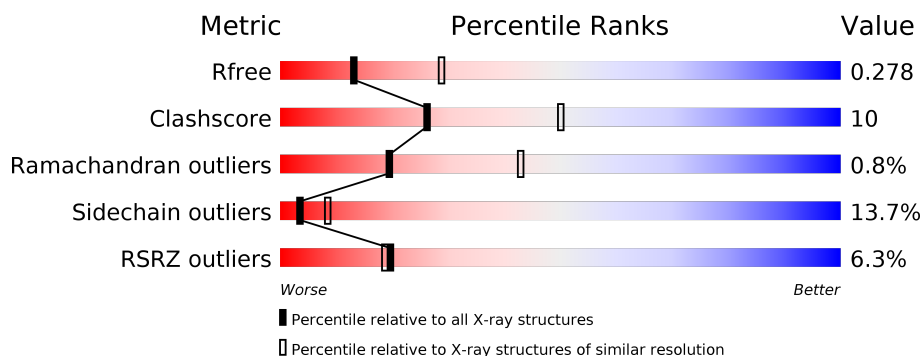
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
$R_{free}$	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	93	<div> <div>3%</div> <div> <div></div> <div>61%</div> <div>20%</div> <div>..</div> <div>16%</div> </div> </div>
1	E	93	<div> <div>13%</div> <div> <div></div> <div>58%</div> <div>12%</div> <div>.</div> <div>27%</div> </div> </div>
2	B	67	<div> <div>3%</div> <div> <div></div> <div>52%</div> <div>31%</div> <div>.</div> <div>15%</div> </div> </div>
2	F	67	<div> <div>6%</div> <div> <div></div> <div>51%</div> <div>37%</div> <div>.</div> <div>9%</div> </div> </div>
3	C	17	<div> <div>12%</div> <div> <div></div> <div>65%</div> <div>24%</div> </div> </div>
3	G	17	<div> <div>41%</div> <div> <div></div> <div>41%</div> <div>18%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
4	D	17	 35% 47% 18%
4	H	17	 24% 41% 35%

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3668 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 Homeodomain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	78	Total	C	N	O	S	0	0	0
			664	414	140	109	1			
1	E	68	Total	C	N	O	S	0	0	0
			580	362	118	99	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	169	MET	-	INITIATING METHIONINE	UNP Q9VD99
E	169	MET	-	INITIATING METHIONINE	UNP Q9VD99

- Molecule 2 is a protein called Homeobox protein aristaless.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	57	Total	C	N	O	S	0	0	0
			504	322	96	85	1			
2	F	61	Total	C	N	O	S	0	0	0
			538	343	102	92	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	17	Total	C	N	O	P	0	0	0
			349	168	63	102	16			
3	G	17	Total	C	N	O	P	0	0	0
			349	168	63	102	16			

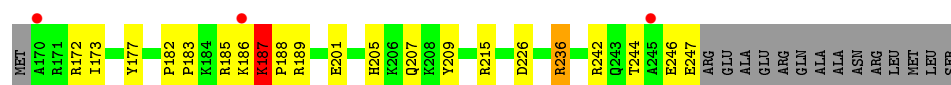
- Molecule 4 is a DNA chain called 5'-D(\*CP\*CP\*GP\*CP\*AP\*AP\*TP\*TP\*AP\*AP\*TP\*TP\*AP\*AP\*GP\*CP\*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	17	Total 342	C 165	N 63	O 98	P 16	0	0	0
4	H	17	Total 342	C 165	N 63	O 98	P 16	0	0	0

### 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 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: Homeodomain-containing protein



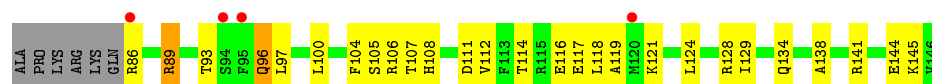
- Molecule 1: Homeodomain-containing protein



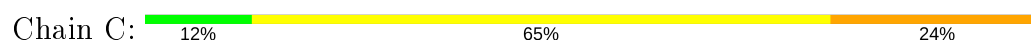
- Molecule 2: Homeobox protein aristaless



- Molecule 2: Homeobox protein aristaless

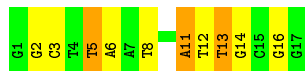


- Molecule 3: 5'-D(\*GP\*GP\*CP\*TP\*TP\*AP\*AP\*TP\*TP\*AP\*AP\*TP\*TP\*GP\*CP\*GP\*G)-3'

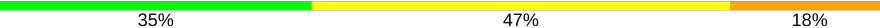


- Molecule 3: 5'-D(\*GP\*GP\*CP\*TP\*TP\*AP\*AP\*TP\*TP\*AP\*AP\*TP\*TP\*GP\*CP\*GP\*G)-3',

Chain G: 




- Molecule 4: 5'-D(\*CP\*CP\*GP\*CP\*AP\*AP\*TP\*TP\*AP\*AP\*TP\*TP\*AP\*AP\*GP\*CP\*C)-3',

Chain D: 



- Molecule 4: 5'-D(\*CP\*CP\*GP\*CP\*AP\*AP\*TP\*TP\*AP\*AP\*TP\*TP\*AP\*AP\*GP\*CP\*C)-3',

Chain H: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.93Å 85.40Å 110.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.65 – 2.70 19.65 – 2.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.65-2.70) 100.0 (19.65-2.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.01 (at 2.71Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.228 , 0.283 0.227 , 0.278	Depositor DCC
$R_{free}$ test set	1943 reflections (10.23%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	64.0	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 34.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3668	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.00% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical 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	0.58	0/678	0.67	0/906
1	E	0.55	0/592	0.65	0/793
2	B	0.60	0/516	0.70	0/691
2	F	0.58	0/550	0.74	0/736
3	C	1.13	2/391 (0.5%)	1.86	11/603 (1.8%)
3	G	1.02	0/391	1.76	10/603 (1.7%)
4	D	1.06	0/383	2.07	23/588 (3.9%)
4	H	1.02	0/383	2.19	16/588 (2.7%)
All	All	0.81	2/3884 (0.1%)	1.40	60/5508 (1.1%)

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

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	6	DA	C3'-O3'	-5.19	1.37	1.44
3	C	14	DG	C3'-O3'	-5.04	1.37	1.44

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	11	DT	O4'-C1'-N1	13.27	117.29	108.00
4	H	7	DT	O4'-C4'-C3'	-11.90	98.86	106.00
4	D	7	DT	C4'-C3'-C2'	-11.78	92.50	103.10
3	G	11	DA	O4'-C4'-C3'	-11.07	99.36	106.00
4	D	7	DT	O4'-C4'-C3'	-10.70	99.58	106.00
3	C	6	DA	O4'-C1'-N9	-10.59	100.59	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	7	DT	C1'-O4'-C4'	-10.17	99.93	110.10
4	H	10	DA	O4'-C1'-N9	10.13	115.09	108.00
4	H	7	DT	C4'-C3'-C2'	-9.95	94.15	103.10
4	H	11	DT	C1'-O4'-C4'	-9.40	100.70	110.10
4	D	14	DA	O4'-C1'-N9	-9.36	101.45	108.00
3	C	14	DG	O4'-C1'-N9	9.35	114.55	108.00
4	D	12	DT	O4'-C1'-N1	-8.42	102.10	108.00
4	H	7	DT	O4'-C1'-N1	8.34	113.84	108.00
4	D	11	DT	C1'-O4'-C4'	-8.05	102.05	110.10
3	C	13	DT	O4'-C1'-N1	-7.98	102.41	108.00
3	G	6	DA	O4'-C1'-N9	-7.90	102.47	108.00
4	D	7	DT	C1'-O4'-C4'	-7.81	102.29	110.10
4	D	11	DT	C4-C5-C7	7.75	123.65	119.00
3	C	10	DA	P-O3'-C3'	7.69	128.93	119.70
4	H	15	DG	O4'-C1'-N9	7.50	113.25	108.00
4	H	1	DC	O4'-C1'-N1	7.40	113.18	108.00
4	D	9	DA	O4'-C1'-N9	-7.39	102.83	108.00
4	D	3	DG	O4'-C1'-N9	-7.34	102.86	108.00
4	D	7	DT	O5'-P-OP1	7.29	119.45	110.70
4	H	10	DA	C1'-O4'-C4'	-6.99	103.11	110.10
3	C	1	DG	O4'-C4'-C3'	-6.95	101.72	104.50
4	H	14	DA	O4'-C1'-N9	-6.71	103.30	108.00
4	H	10	DA	O4'-C1'-C2'	-6.55	100.66	105.90
3	G	5	DT	C4-C5-C7	6.52	122.91	119.00
3	C	10	DA	O4'-C1'-N9	-6.39	103.53	108.00
3	G	5	DT	C6-C5-C7	-6.29	119.13	122.90
4	H	11	DT	O4'-C1'-C2'	-6.28	100.88	105.90
3	G	5	DT	O4'-C1'-N1	6.17	112.32	108.00
3	G	8	DT	N3-C2-O2	-6.01	118.69	122.30
4	H	11	DT	C3'-C2'-C1'	-6.00	95.31	102.50
4	D	10	DA	O4'-C1'-N9	5.98	112.18	108.00
4	D	10	DA	P-O5'-C5'	-5.97	111.35	120.90
4	D	7	DT	C5'-C4'-O4'	5.92	120.55	109.30
3	G	11	DA	P-O3'-C3'	5.89	126.77	119.70
4	D	11	DT	O4'-C4'-C3'	-5.82	102.17	104.50
3	G	13	DT	O4'-C1'-N1	-5.82	103.93	108.00
4	D	11	DT	O4'-C1'-N1	5.76	112.03	108.00
3	C	5	DT	O4'-C1'-N1	5.70	111.99	108.00
4	D	7	DT	O4'-C1'-N1	5.45	111.82	108.00
4	D	15	DG	O4'-C1'-N9	5.43	111.80	108.00
4	D	11	DT	C6-C5-C7	-5.40	119.66	122.90
3	C	17	DG	C1'-O4'-C4'	-5.39	104.71	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	12	DT	N3-C4-O4	5.34	123.10	119.90
4	D	10	DA	C5'-C4'-C3'	-5.33	104.50	114.10
4	D	15	DG	C3'-C2'-C1'	-5.33	96.10	102.50
3	G	16	DG	O4'-C1'-N9	5.32	111.72	108.00
4	D	7	DT	N3-C2-O2	-5.23	119.16	122.30
3	C	4	DT	N3-C4-O4	5.20	123.02	119.90
4	H	9	DA	P-O3'-C3'	5.15	125.89	119.70
3	G	11	DA	O4'-C1'-N9	5.11	111.58	108.00
4	D	9	DA	O4'-C4'-C3'	-5.09	102.46	104.50
4	H	5	DA	C2-N3-C4	5.09	113.14	110.60
3	C	17	DG	O4'-C1'-N9	5.03	111.52	108.00
3	C	3	DC	O4'-C1'-N1	-5.02	104.48	108.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	187	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	664	0	689	16	0
1	E	580	0	584	12	0
2	B	504	0	503	12	0
2	F	538	0	539	13	0
3	C	349	0	195	7	0
3	G	349	0	195	6	0
4	D	342	0	193	11	0
4	H	342	0	193	14	0
All	All	3668	0	3091	70	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 10.

All (70) 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:A:186:LYS:HA	1:A:187:LYS:HB3	1.37	1.04
2:B:93:THR:OG1	2:B:96:GLN:HG2	1.73	0.88
1:E:201:GLU:OE2	1:E:236:ARG:NH1	2.15	0.80
2:F:89:ARG:NH2	3:G:5:DT:O2	2.15	0.79
1:A:242:ARG:O	1:A:246:GLU:HB2	1.82	0.79
1:E:177:TYR:HB2	4:H:7:DT:H5'	1.65	0.79
4:D:10:DA:OP1	2:F:145:LYS:HG3	1.82	0.78
1:A:177:TYR:HB3	4:D:7:DT:H5'	1.67	0.76
1:E:215:ARG:NH1	4:H:2:DC:OP1	2.18	0.75
1:A:186:LYS:HA	1:A:187:LYS:CB	2.16	0.67
1:A:177:TYR:CB	4:D:7:DT:H5'	2.24	0.66
1:E:177:TYR:CB	4:H:7:DT:H5'	2.29	0.62
2:F:119:ALA:HB2	2:F:129:ILE:HD12	1.83	0.61
2:F:100:LEU:HD21	2:F:129:ILE:HG12	1.84	0.60
3:C:16:DG:H2''	3:C:17:DG:OP2	2.03	0.58
3:G:11:DA:H2''	3:G:12:DT:O5'	2.02	0.58
3:C:12:DT:H2'	3:C:13:DT:H72	1.85	0.57
1:E:180:ARG:HG2	3:G:14:DG:H5'	1.86	0.57
1:A:215:ARG:NH1	4:D:2:DC:OP1	2.29	0.56
1:E:176:PRO:HB2	1:E:178:GLN:HE21	1.70	0.56
2:B:109:TYR:CE1	4:D:7:DT:H5''	2.43	0.54
1:A:201:GLU:OE1	1:A:236:ARG:HD3	2.08	0.54
2:F:138:ALA:HA	2:F:141:ARG:NH1	2.22	0.54
3:G:11:DA:C2'	3:G:12:DT:O5'	2.56	0.54
3:C:9:DT:H2''	3:C:10:DA:C8	2.44	0.53
2:B:95:PHE:O	2:B:98:GLU:HB3	2.09	0.52
2:B:104:PHE:CZ	2:B:108:HIS:HD2	2.26	0.52
3:C:2:DG:H1'	3:C:3:DC:H5'	1.90	0.52
4:H:11:DT:OP2	4:H:11:DT:H2'	2.11	0.51
1:E:215:ARG:CZ	1:E:226:ASP:OD1	2.60	0.50
2:B:140:TRP:O	2:F:107:THR:HG22	2.12	0.49
3:G:2:DG:H2''	3:G:3:DC:C6	2.47	0.49
1:A:187:LYS:H	1:A:188:PRO:HD2	1.77	0.49
2:B:111:ASP:OD2	2:B:114:THR:OG1	2.30	0.49
2:F:111:ASP:OD1	2:F:114:THR:OG1	2.24	0.49
4:H:5:DA:H1'	4:H:6:DA:H5'	1.95	0.49
1:A:187:LYS:HG2	1:A:188:PRO:N	2.28	0.48
2:B:109:TYR:HE1	4:D:7:DT:H5''	1.78	0.48
3:C:12:DT:H2'	3:C:13:DT:C7	2.44	0.48
4:H:1:DC:H5''	4:H:1:DC:H6	1.80	0.47
4:H:8:DT:H2''	4:H:9:DA:C8	2.50	0.46
1:A:177:TYR:CG	4:D:7:DT:H4'	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:ILE:HG13	4:D:6:DA:OP1	2.16	0.46
3:G:12:DT:H2'	3:G:13:DT:H71	1.98	0.46
2:F:93:THR:OG1	2:F:96:GLN:HG2	2.15	0.45
1:E:189:ARG:HH12	4:H:9:DA:H2	1.65	0.45
4:D:10:DA:H2''	4:D:11:DT:O4'	2.17	0.45
1:A:185:ARG:HD3	2:F:144:GLU:O	2.17	0.45
3:C:11:DA:H2'	3:C:12:DT:C6	2.53	0.44
1:E:215:ARG:NH2	1:E:226:ASP:OD2	2.52	0.43
2:B:141:ARG:NH2	4:D:8:DT:OP1	2.52	0.43
1:A:172:ARG:HD2	1:A:172:ARG:H	1.83	0.43
1:E:225:THR:H	1:E:228:GLN:HE21	1.67	0.43
4:D:10:DA:C2	4:D:11:DT:C2	3.07	0.43
1:A:172:ARG:HA	1:A:172:ARG:NE	2.34	0.42
4:H:2:DC:H2'	4:H:3:DG:C8	2.54	0.42
1:A:182:PRO:HA	1:A:183:PRO:HD3	1.86	0.42
3:C:15:DC:H2''	3:C:16:DG:C8	2.54	0.42
1:E:177:TYR:HB2	4:H:7:DT:C5'	2.45	0.42
2:B:104:PHE:CZ	2:B:108:HIS:CD2	3.08	0.42
2:F:141:ARG:NH2	4:H:8:DT:OP1	2.53	0.42
2:B:92:PHE:HB3	2:B:96:GLN:HG3	2.01	0.41
2:F:104:PHE:CE1	2:F:108:HIS:ND1	2.89	0.41
2:B:119:ALA:HB1	2:B:124:LEU:O	2.21	0.41
2:F:119:ALA:HB1	2:F:124:LEU:O	2.20	0.41
4:H:13:DA:H2''	4:H:14:DA:C8	2.56	0.41
1:E:177:TYR:HB2	4:H:7:DT:H4'	2.03	0.41
2:B:96:GLN:HB2	2:B:122:ILE:HD12	2.02	0.41
2:F:134:GLN:HG2	4:H:8:DT:C7	2.51	0.40
1:A:207:GLN:HG2	1:A:209:TYR:O	2.21	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	76/93 (82%)	70 (92%)	4 (5%)	2 (3%)	5	13
1	E	64/93 (69%)	62 (97%)	2 (3%)	0	100	100
2	B	55/67 (82%)	52 (94%)	3 (6%)	0	100	100
2	F	59/67 (88%)	55 (93%)	4 (7%)	0	100	100
All	All	254/320 (79%)	239 (94%)	13 (5%)	2 (1%)	19	43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	187	LYS
1	A	244	THR

### 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	67/79 (85%)	61 (91%)	6 (9%)	9	22
1	E	59/79 (75%)	53 (90%)	6 (10%)	7	17
2	B	52/61 (85%)	44 (85%)	8 (15%)	2	7
2	F	56/61 (92%)	44 (79%)	12 (21%)	1	3
All	All	234/280 (84%)	202 (86%)	32 (14%)	3	8

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	187	LYS
1	A	189	ARG
1	A	205	HIS
1	A	226	ASP
1	A	236	ARG
1	A	247	GLU
2	B	89	ARG
2	B	97	LEU
2	B	107	THR

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Mol	Chain	Res	Type
2	B	112	VAL
2	B	118	LEU
2	B	121	LYS
2	B	126	GLU
2	B	140	TRP
1	E	178	GLN
1	E	194	ARG
1	E	208	LYS
1	E	215	ARG
1	E	236	ARG
1	E	242	ARG
2	F	86	ARG
2	F	89	ARG
2	F	96	GLN
2	F	97	LEU
2	F	105	SER
2	F	106	ARG
2	F	112	VAL
2	F	116	GLU
2	F	117	GLU
2	F	118	LEU
2	F	121	LYS
2	F	128	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	108	HIS
2	B	130	GLN
2	B	134	GLN
1	E	178	GLN
1	E	228	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 [i](#)

There are no carbohydrates 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, 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	78/93 (83%)	0.27	3 (3%) 40 39	35, 54, 76, 80	0
1	E	68/93 (73%)	0.61	12 (17%) 1 1	35, 56, 109, 110	0
2	B	57/67 (85%)	-0.08	2 (3%) 44 44	32, 53, 65, 78	0
2	F	61/67 (91%)	0.26	4 (6%) 18 16	48, 61, 82, 90	0
3	C	17/17 (100%)	-0.51	0 100 100	41, 46, 62, 66	0
3	G	17/17 (100%)	-0.12	0 100 100	43, 57, 72, 72	0
4	D	17/17 (100%)	-0.37	0 100 100	36, 52, 67, 68	0
4	H	17/17 (100%)	-0.20	0 100 100	43, 52, 69, 75	0
All	All	332/388 (85%)	0.16	21 (6%) 20 19	32, 56, 81, 110	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	190	THR	6.8
1	E	175	HIS	6.0
1	E	181	THR	5.9
2	B	86	ARG	4.6
1	E	182	PRO	4.4
2	F	86	ARG	4.4
1	A	170	ALA	4.2
1	E	189	ARG	4.1
1	E	245	ALA	3.4
1	E	176	PRO	2.9
1	E	178	GLN	2.9
2	F	95	PHE	2.7
1	E	247	GLU	2.7
1	A	186	LYS	2.6
1	E	177	TYR	2.6
1	E	183	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	120	MET	2.4
1	A	245	ALA	2.2
1	E	179	ASN	2.2
2	F	120	MET	2.1
2	F	94	SER	2.1

## 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 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.