



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:04 PM UTC

PDB ID : 8IPF / pdb\_00008ipf  
Title : Crystal structure of an ankyrin protein that can specifically recognize a parallel G-quadruplex  
Authors : Ngo, K.H.; Liew, C.W.; Heddi, B.; Phan, A.T.  
Deposited on : 2023-03-14  
Resolution : 2.01 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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

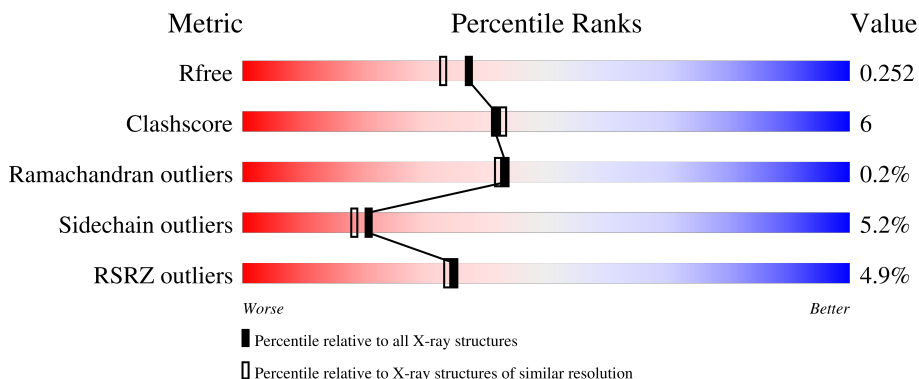
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.01 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.



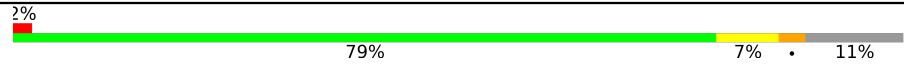


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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  $\geq 3$ , 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	136	 3% 76% 11% • 10%
1	B	136	 7% 73% 15% • 11%
1	C	136	 4% 75% 13% • 11%
1	D	136	 3% 68% 20% • 11%
1	E	136	 3% 76% 12% • 11%

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Mol	Chain	Length	Quality of chain
1	F	136	
1	G	136	
1	H	136	

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 7918 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 DARPin (2E4).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	122	Total 936	C 585	N 168	O 180	S 3	0	0	0
1	B	121	Total 934	C 584	N 168	O 179	S 3	0	1	0
1	C	121	Total 943	C 588	N 171	O 181	S 3	0	2	0
1	D	121	Total 930	C 581	N 166	O 180	S 3	0	1	0
1	E	121	Total 927	C 579	N 166	O 179	S 3	0	0	0
1	F	121	Total 938	C 585	N 170	O 180	S 3	0	1	0
1	G	121	Total 953	C 598	N 172	O 179	S 4	0	4	0
1	H	121	Total 934	C 584	N 168	O 179	S 3	0	1	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	58	Total 58	O 58	0	0
2	B	54	Total 54	O 54	0	0
2	C	53	Total 53	O 53	0	0
2	D	59	Total 59	O 59	0	0
2	E	53	Total 53	O 53	0	0
2	F	62	Total 62	O 62	0	0

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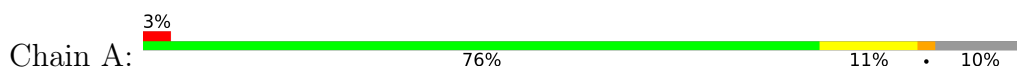
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
2	G	37	Total	O	0	0
			37	37		
2	H	47	Total	O	0	0
			47	47		

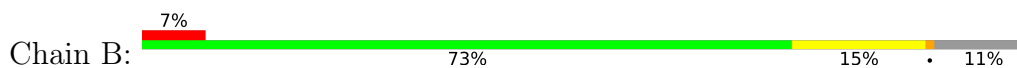
### 3 Residue-property plots [i](#)

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.

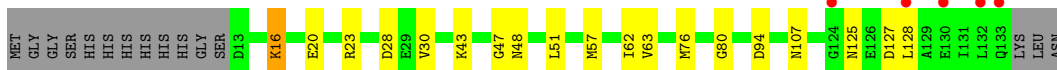
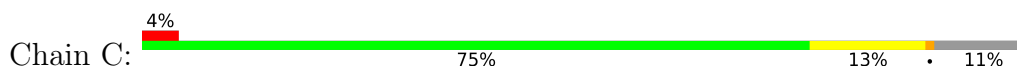
- Molecule 1: DARPin (2E4)



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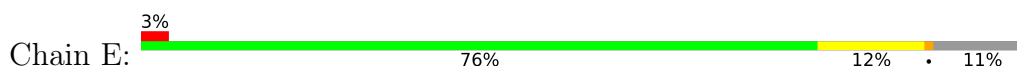


- Molecule 1: DARPin (2E4)



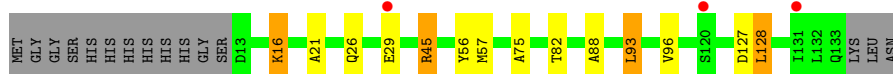
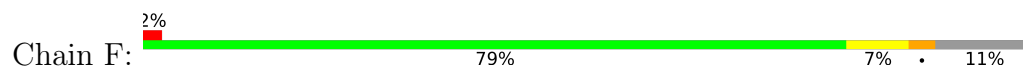
LYS  
LEU  
ASN

- Molecule 1: DARPin (2E4)

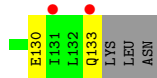




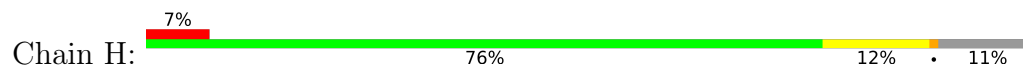
- Molecule 1: DARPin (2E4)



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- Molecule 1: DARPin (2E4)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	64.30Å 59.07Å 127.50Å 90.00° 100.56° 90.00°	Depositor
Resolution (Å)	49.25 – 2.01 49.25 – 2.01	Depositor EDS
% Data completeness (in resolution range)	99.1 (49.25-2.01) 99.1 (49.25-2.01)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.62 (at 2.01Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, $R_{free}$	0.226 , 0.249 0.228 , 0.252	Depositor DCC
$R_{free}$ test set	3127 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	29.0	Xtrriage
Anisotropy	1.191	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 44.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.55$ , $\langle L^2 \rangle = 0.39$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7918	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	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 50.40 % 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 6.5006e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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	0.85	0/950	1.15	0/1283
1	B	0.82	0/952	1.10	2/1287 (0.2%)
1	C	0.90	2/960 (0.2%)	1.20	2/1297 (0.2%)
1	D	0.94	2/947 (0.2%)	1.22	2/1280 (0.2%)
1	E	0.85	0/941	1.18	0/1272
1	F	0.79	0/952	1.07	1/1286 (0.1%)
1	G	1.00	4/980 (0.4%)	1.38	10/1322 (0.8%)
1	H	0.79	0/952	1.14	2/1287 (0.2%)
All	All	0.87	8/7634 (0.1%)	1.18	19/10314 (0.2%)

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	F	0	2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	31[A]	ARG	C-O	6.83	1.32	1.24
1	G	31[B]	ARG	C-O	6.83	1.32	1.24
1	D	120[A]	SER	C-O	6.67	1.31	1.24
1	D	120[B]	SER	C-O	6.67	1.31	1.24
1	G	43[A]	LYS	C-O	5.73	1.31	1.24
1	G	43[B]	LYS	C-O	5.73	1.31	1.24
1	C	107[A]	ASN	C-O	5.61	1.31	1.24
1	C	107[B]	ASN	C-O	5.61	1.31	1.24

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	107[A]	ASN	CA-C-O	8.03	129.14	119.43
1	C	107[B]	ASN	CA-C-O	8.03	129.14	119.43
1	G	43[A]	LYS	CA-C-O	7.64	128.95	120.70
1	G	43[B]	LYS	CA-C-O	7.64	128.95	120.70
1	D	120[A]	SER	CA-C-O	7.17	128.38	120.63
1	D	120[B]	SER	CA-C-O	7.17	128.38	120.63
1	G	31[A]	ARG	CA-C-O	7.07	128.05	120.55
1	G	31[B]	ARG	CA-C-O	7.07	128.05	120.55
1	G	78[A]	HIS	CA-C-O	6.79	127.70	119.49
1	G	78[B]	HIS	CA-C-O	6.79	127.70	119.49
1	B	78[A]	HIS	CA-C-O	6.40	127.25	119.38
1	B	78[B]	HIS	CA-C-O	6.40	127.25	119.38
1	G	34[A]	MET	CA-C-O	6.36	127.50	120.63
1	G	34[B]	MET	CA-C-O	6.36	127.50	120.63
1	H	78[A]	HIS	CA-C-O	5.67	126.25	119.10
1	H	78[B]	HIS	CA-C-O	5.67	126.25	119.10
1	F	93	LEU	N-CA-C	5.07	116.50	111.07
1	G	43[A]	LYS	N-CA-C	5.02	117.59	109.40
1	G	43[B]	LYS	N-CA-C	5.02	117.59	109.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	45[A]	ARG	Mainchain
1	F	45[B]	ARG	Mainchain

## 5.2 Too-close contacts [\(i\)](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	936	0	924	11	1
1	B	934	0	918	10	0
1	C	943	0	929	14	0
1	D	930	0	916	16	0
1	E	927	0	911	10	0
1	F	938	0	923	14	0
1	G	953	0	953	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	934	0	918	7	1
2	A	58	0	0	4	0
2	B	54	0	0	0	0
2	C	53	0	0	5	0
2	D	59	0	0	1	0
2	E	53	0	0	1	0
2	F	62	0	0	4	0
2	G	37	0	0	0	0
2	H	47	0	0	2	0
All	All	7918	0	7392	95	1

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 6.

All (95) 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:C:94:ASP:HB3	2:C:231:HOH:O	1.70	0.90
1:F:57:MET:HE3	2:F:216:HOH:O	1.75	0.86
1:E:56:TYR:CD1	1:E:57:MET:HE2	2.13	0.83
1:A:81:ARG:HD3	2:A:235:HOH:O	1.79	0.82
1:F:56:TYR:CD1	1:F:57:MET:HE2	2.17	0.79
1:E:56:TYR:HD1	1:E:57:MET:HE2	1.45	0.78
1:F:56:TYR:HD1	1:F:57:MET:HE2	1.54	0.71
1:B:45:ARG:HD3	2:H:222:HOH:O	1.94	0.67
1:H:56:TYR:CD2	1:H:57:MET:HE2	2.30	0.67
1:D:93:LEU:HG	1:D:128:LEU:HD13	1.77	0.67
1:H:57:MET:HE3	2:H:201:HOH:O	1.95	0.66
1:H:56:TYR:HD2	1:H:57:MET:HE2	1.63	0.63
1:A:81:ARG:CD	2:A:235:HOH:O	2.43	0.63
1:E:130:GLU:HA	1:E:133:GLN:HG2	1.82	0.61
1:A:26:GLN:HB3	1:A:29:GLU:HB2	1.82	0.61
1:C:94:ASP:CA	2:C:231:HOH:O	2.49	0.60
1:C:94:ASP:CB	2:C:231:HOH:O	2.39	0.60
1:F:45[B]:ARG:NH2	2:F:203:HOH:O	2.35	0.59
1:C:94:ASP:HA	2:C:231:HOH:O	2.02	0.59
1:D:130:GLU:HG3	1:E:78:HIS:CE1	2.37	0.59
1:C:16:LYS:HE3	1:C:20:GLU:HG3	1.85	0.58
1:F:56:TYR:CE1	1:F:57:MET:HE2	2.38	0.58
1:A:78:HIS:HE1	2:A:251:HOH:O	1.87	0.58
1:C:30:VAL:HG21	1:C:62:ILE:HD13	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:88:ALA:HB1	1:E:128:LEU:HD21	1.88	0.56
1:H:127:ASP:O	1:H:131:ILE:HD12	2.06	0.56
1:G:78[B]:HIS:HD2	1:G:79:TRP:CE2	2.22	0.55
1:E:13:ASP:O	1:E:17:LYS:HG3	2.09	0.53
1:H:84:LEU:HD11	1:H:96:VAL:HG13	1.89	0.53
1:G:44:ASP:CG	1:G:48:ASN:HB2	2.33	0.53
1:D:131:ILE:O	1:D:133:GLN:HG2	2.09	0.53
1:H:88:ALA:HA	1:H:128:LEU:HD21	1.91	0.52
1:E:57:MET:HE3	2:E:208:HOH:O	2.09	0.52
1:D:28:ASP:O	1:D:32:ILE:HG12	2.10	0.51
1:D:78:HIS:HB3	2:D:248:HOH:O	2.10	0.51
1:D:61:GLU:O	1:D:65:VAL:HG13	2.11	0.51
1:F:21:ALA:HA	1:F:26:GLN:HE21	1.76	0.50
1:C:76:MET:HE3	1:C:80:GLY:HA2	1.93	0.50
1:C:43:LYS:HE3	1:C:47:GLY:HA2	1.94	0.49
1:E:43:LYS:HD3	1:E:49:THR:HG22	1.94	0.49
1:A:79:TRP:C	1:A:111:LYS:HG3	2.38	0.49
1:D:42:ALA:O	1:D:50:PRO:HD3	2.12	0.49
1:F:93:LEU:HD21	1:F:127:ASP:HB3	1.95	0.48
1:C:125:ASN:HB3	1:C:128:LEU:HB3	1.95	0.48
1:E:30:VAL:HG21	1:E:62:ILE:HD13	1.95	0.48
1:D:93:LEU:HD12	1:D:93:LEU:H	1.78	0.48
1:D:56:TYR:CD1	1:D:57:MET:HE2	2.49	0.47
1:F:57:MET:CE	2:F:216:HOH:O	2.48	0.47
1:H:100:LEU:HD23	1:H:104:ALA:HB3	1.97	0.47
1:F:16:LYS:HD2	2:F:261:HOH:O	2.15	0.46
1:B:124:GLY:O	1:B:126:GLU:N	2.49	0.46
1:D:131:ILE:C	1:D:133:GLN:H	2.23	0.45
1:G:120:SER:OG	1:G:125:ASN:HB2	2.17	0.45
1:F:16:LYS:N	1:F:16:LYS:HD3	2.31	0.45
1:G:68:LYS:HD3	1:G:69:TYR:CE1	2.51	0.45
1:B:127:ASP:N	1:B:127:ASP:OD1	2.50	0.45
1:G:85:HIS:HD2	1:G:119:ILE:HD12	1.82	0.45
1:C:51:LEU:HD11	1:C:63:VAL:HG13	1.98	0.45
1:G:130:GLU:HA	1:G:133:GLN:HG2	1.99	0.45
2:A:219:HOH:O	1:B:78[B]:HIS:CE1	2.70	0.44
1:A:13:ASP:O	1:A:17:LYS:HG2	2.18	0.44
1:A:130:GLU:O	1:A:134:LYS:HG3	2.17	0.43
1:B:25:GLY:HA2	1:B:62:ILE:HG13	2.01	0.43
1:D:22:ALA:HA	1:D:62:ILE:HD13	2.00	0.43
1:F:75:ALA:O	1:F:82:THR:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:94:ASP:OD1	1:B:94:ASP:N	2.51	0.43
1:A:93:LEU:HA	1:A:128:LEU:HD21	1.99	0.43
1:D:131:ILE:C	1:D:133:GLN:N	2.75	0.43
1:F:88:ALA:HA	1:F:128:LEU:HD21	2.01	0.43
1:C:23[B]:ARG:HG3	1:C:57:MET:HE3	2.01	0.43
1:B:97:GLU:O	1:B:101:LYS:HG3	2.19	0.42
1:G:42:ALA:O	1:G:50:PRO:HD3	2.19	0.42
1:A:84:LEU:HD11	1:A:96:VAL:HG13	2.02	0.42
1:C:43:LYS:HA	1:C:48:ASN:O	2.20	0.42
1:B:16:LYS:O	1:B:20:GLU:HG3	2.19	0.42
1:G:65:VAL:O	1:G:69:TYR:HD1	2.03	0.42
1:G:78[B]:HIS:HD2	1:G:79:TRP:NE1	2.18	0.42
1:D:14:LEU:HB3	1:D:38:ALA:HB2	2.02	0.41
1:D:43:LYS:HA	1:D:48:ASN:O	2.20	0.41
1:B:75:ALA:O	1:B:82:THR:HA	2.20	0.41
1:D:44:ASP:CG	1:D:48:ASN:HB2	2.45	0.41
1:A:130:GLU:HG3	1:A:134:LYS:HE3	2.02	0.41
1:G:44:ASP:OD2	1:G:48:ASN:HB2	2.21	0.41
1:B:14:LEU:HD22	1:B:14:LEU:HA	1.96	0.41
1:F:96:VAL:HG21	1:F:128:LEU:HD11	2.02	0.41
1:A:16:LYS:HE2	1:A:16:LYS:HB2	1.57	0.41
1:C:62:ILE:HG13	2:C:214:HOH:O	2.20	0.41
1:D:27:ASP:O	1:D:31:ARG:HG3	2.20	0.41
1:E:42:ALA:O	1:E:50:PRO:HD3	2.21	0.41
1:G:22:ALA:O	1:G:59:HIS:HD2	2.04	0.41
1:G:130:GLU:HA	1:G:133:GLN:HE21	1.86	0.40
1:G:27:ASP:O	1:G:31[B]:ARG:HG3	2.22	0.40
1:F:128:LEU:HD12	1:F:128:LEU:HA	1.82	0.40
1:G:94:ASP:OD1	1:G:94:ASP:N	2.54	0.40
1:C:127:ASP:N	1:C:127:ASP:OD1	2.54	0.40

All (1) 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)
1:A:29:GLU:OE1	1:H:69:TYR:OH[2_546]	2.06	0.14

## 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	120/136 (88%)	119 (99%)	1 (1%)	0	100	100
1	B	120/136 (88%)	119 (99%)	0	1 (1%)	16	11
1	C	121/136 (89%)	118 (98%)	3 (2%)	0	100	100
1	D	120/136 (88%)	119 (99%)	1 (1%)	0	100	100
1	E	119/136 (88%)	118 (99%)	1 (1%)	0	100	100
1	F	120/136 (88%)	119 (99%)	1 (1%)	0	100	100
1	G	123/136 (90%)	122 (99%)	0	1 (1%)	16	11
1	H	120/136 (88%)	120 (100%)	0	0	100	100
All	All	963/1088 (88%)	954 (99%)	7 (1%)	2 (0%)	43	42

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	125	ASN
1	G	125	ASN

### 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	95/106 (90%)	89 (94%)	6 (6%)	16	13
1	B	95/106 (90%)	88 (93%)	7 (7%)	13	9
1	C	96/106 (91%)	94 (98%)	2 (2%)	47	52

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	95/106 (90%)	90 (95%)	5 (5%)	20	18
1	E	94/106 (89%)	90 (96%)	4 (4%)	26	25
1	F	95/106 (90%)	92 (97%)	3 (3%)	34	35
1	G	98/106 (92%)	93 (95%)	5 (5%)	21	19
1	H	95/106 (90%)	88 (93%)	7 (7%)	13	9
All	All	763/848 (90%)	724 (95%)	39 (5%)	21	19

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LYS
1	A	26	GLN
1	A	27	ASP
1	A	76	MET
1	A	130	GLU
1	A	131	ILE
1	B	13	ASP
1	B	14	LEU
1	B	26	GLN
1	B	27	ASP
1	B	32	ILE
1	B	33	LEU
1	B	131	ILE
1	C	16	LYS
1	C	28	ASP
1	D	16	LYS
1	D	106	VAL
1	D	127	ASP
1	D	131	ILE
1	D	132	LEU
1	E	16	LYS
1	E	64	GLU
1	E	126	GLU
1	E	130	GLU
1	F	16	LYS
1	F	29	GLU
1	F	128	LEU
1	G	13	ASP
1	G	93	LEU
1	G	111	LYS

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Mol	Chain	Res	Type
1	G	114	LYS
1	G	126	GLU
1	H	31	ARG
1	H	93	LEU
1	H	94	ASP
1	H	101	LYS
1	H	111	LYS
1	H	130	GLU
1	H	131	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	78	HIS
1	B	123	ASN
1	C	92	HIS
1	C	102	HIS
1	D	48	ASN
1	D	102	HIS
1	D	109	GLN
1	D	133	GLN
1	E	48	ASN
1	F	26	GLN
1	F	48	ASN
1	G	48	ASN
1	G	123	ASN
1	G	133	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, 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	122/136 (89%)	0.49	4 (3%) 49 48	29, 38, 50, 68	0
1	B	121/136 (88%)	0.70	9 (7%) 20 19	21, 40, 70, 81	1 (0%)
1	C	121/136 (88%)	0.68	5 (4%) 41 40	21, 39, 60, 73	2 (1%)
1	D	121/136 (88%)	0.62	4 (3%) 49 48	31, 39, 58, 69	1 (0%)
1	E	121/136 (88%)	0.64	4 (3%) 49 48	31, 41, 62, 71	0
1	F	121/136 (88%)	0.59	3 (2%) 58 58	21, 39, 56, 71	1 (0%)
1	G	121/136 (88%)	0.84	8 (6%) 24 23	23, 43, 69, 79	4 (3%)
1	H	121/136 (88%)	0.81	10 (8%) 17 16	22, 43, 62, 70	1 (0%)
All	All	969/1088 (89%)	0.67	47 (4%) 35 34	21, 40, 62, 81	10 (1%)

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	127	ASP	3.7
1	H	93	LEU	3.6
1	A	13	ASP	3.3
1	E	131	ILE	3.1
1	F	29	GLU	2.8
1	C	133	GLN	2.8
1	C	128	LEU	2.7
1	B	26	GLN	2.7
1	F	120	SER	2.6
1	H	133	GLN	2.6
1	A	27	ASP	2.6
1	A	112	PHE	2.6
1	E	89	SER	2.6
1	D	132	LEU	2.5
1	B	93	LEU	2.5
1	G	131	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	125	ASN	2.4
1	D	125	ASN	2.4
1	H	129	ALA	2.4
1	F	131	ILE	2.4
1	B	130	GLU	2.4
1	G	94	ASP	2.3
1	H	98	VAL	2.3
1	C	132	LEU	2.3
1	A	134	LYS	2.3
1	B	124	GLY	2.3
1	B	112	PHE	2.2
1	H	127	ASP	2.2
1	G	124	GLY	2.2
1	D	93	LEU	2.2
1	C	124	GLY	2.2
1	H	15	GLY	2.1
1	H	35	ALA	2.1
1	G	73	VAL	2.1
1	B	129	ALA	2.1
1	H	106	VAL	2.1
1	C	130	GLU	2.0
1	H	41	ASN	2.0
1	B	13	ASP	2.0
1	B	122	ASP	2.0
1	G	133	GLN	2.0
1	E	124	GLY	2.0
1	D	65	VAL	2.0
1	G	89	SER	2.0
1	H	88	ALA	2.0
1	G	125	ASN	2.0
1	E	128	LEU	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

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.