



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:38 AM GMT

PDB ID : 2NNN  
Title : Crystal structure of probable transcriptional regulator from *Pseudomonas aeruginosa*  
Authors : Chang, C.; Evdokimova, E.; Altamentova, S.; Savchenko, A.; Edwards, A.M.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)  
Deposited on : 2006-10-24  
Resolution : 2.40 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  
<http://wwpdb.org/validation/2016/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  
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

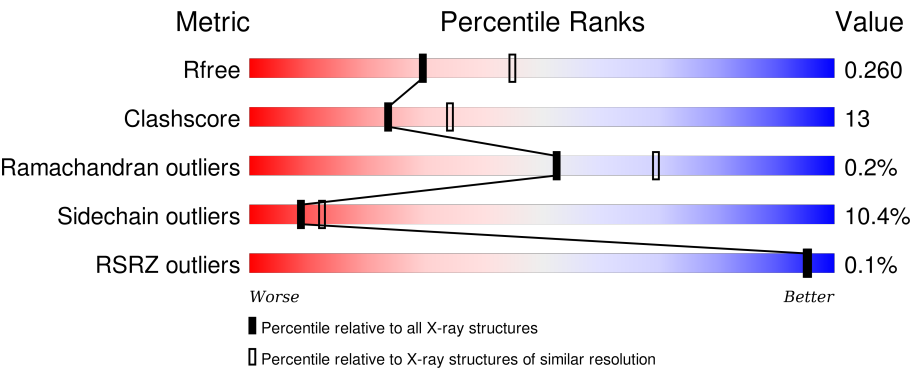
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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 <sub>free</sub>	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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. 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	140	
1	B	140	
1	C	140	
1	D	140	
1	E	140	

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Mol	Chain	Length	Quality of chain
1	F	140	<div><div></div><div>69%22%5%</div></div>
1	G	140	<div>%<div><div></div><div>59%31%6%</div></div></div>
1	H	140	<div><div></div><div>75%19%5%</div></div>
1	I	140	<div><div></div><div>72%21%5%</div></div>
1	J	140	<div><div></div><div>71%21%5%</div></div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10506 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 Probable transcriptional regulator.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	134	Total	C	N	O	S	Se	0	0	0
			1009	624	194	189	1	1			
1	B	134	Total	C	N	O	S	Se	0	0	0
			1009	624	194	189	1	1			
1	C	132	Total	C	N	O	S	Se	0	0	0
			995	616	192	185	1	1			
1	D	133	Total	C	N	O	S	Se	0	0	0
			1004	621	193	188	1	1			
1	E	132	Total	C	N	O	S	Se	0	0	0
			996	617	192	185	1	1			
1	F	133	Total	C	N	O	S	Se	0	0	0
			1004	621	193	188	1	1			
1	G	132	Total	C	N	O	S	Se	0	0	0
			995	616	192	185	1	1			
1	H	133	Total	C	N	O	S	Se	0	0	0
			1004	621	193	188	1	1			
1	I	133	Total	C	N	O	S	Se	0	0	0
			1004	621	193	188	1	1			
1	J	133	Total	C	N	O	S	Se	0	0	0
			1004	621	193	188	1	1			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
A	64	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
B	64	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
C	1	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
C	64	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
D	1	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
D	64	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
E	1	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8

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Chain	Residue	Modelled	Actual	Comment	Reference
E	64	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
F	1	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
F	64	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
G	1	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
G	64	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
H	1	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
H	64	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
I	1	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
I	64	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
J	1	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8
J	64	MSE	MET	MODIFIED RESIDUE	UNP Q9I3B8

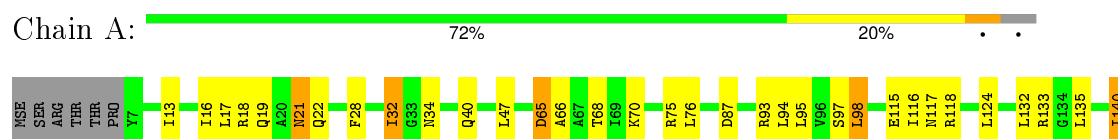
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	67	Total O 67 67	0	0
2	B	62	Total O 62 62	0	0
2	C	36	Total O 36 36	0	0
2	D	63	Total O 63 63	0	0
2	E	45	Total O 45 45	0	0
2	F	38	Total O 38 38	0	0
2	G	42	Total O 42 42	0	0
2	H	51	Total O 51 51	0	0
2	I	33	Total O 33 33	0	0
2	J	45	Total O 45 45	0	0

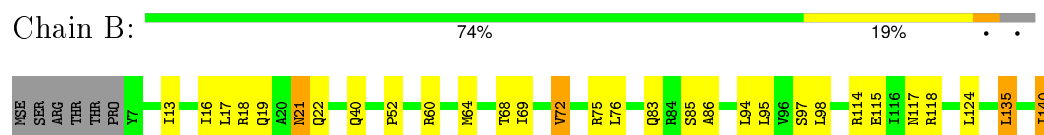
### 3 Residue-property plots

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.

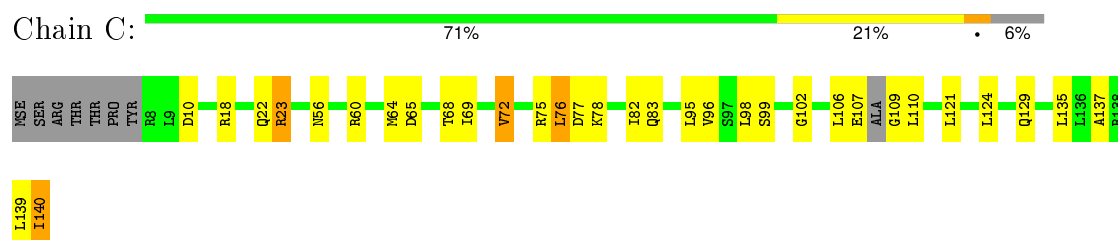
- Molecule 1: Probable transcriptional regulator



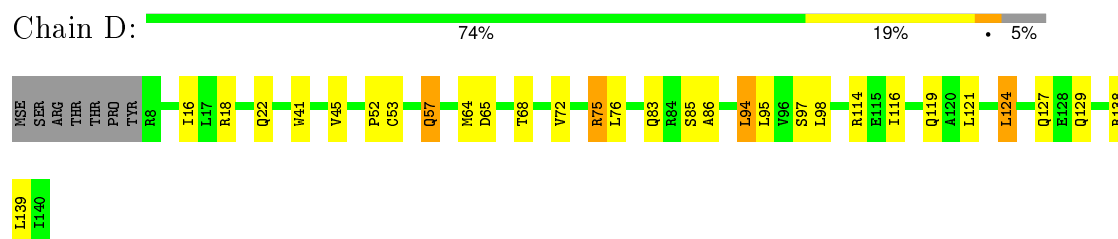
- Molecule 1: Probable transcriptional regulator



- Molecule 1: Probable transcriptional regulator

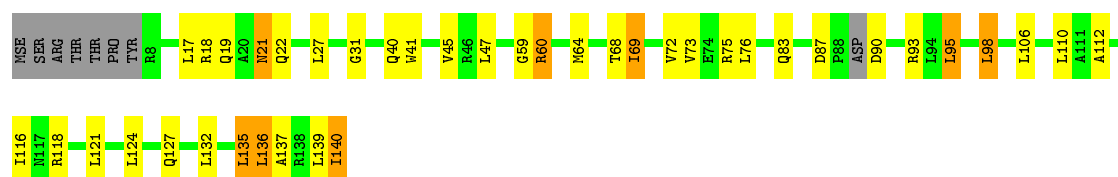


- Molecule 1: Probable transcriptional regulator



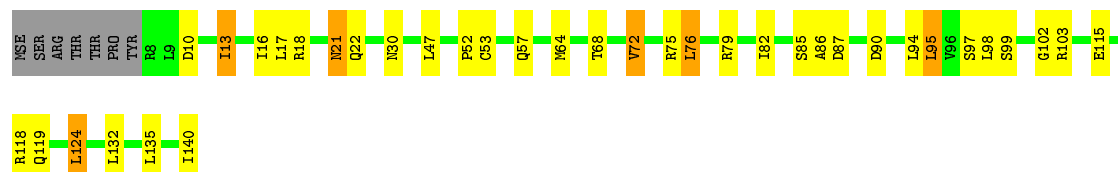
- Molecule 1: Probable transcriptional regulator





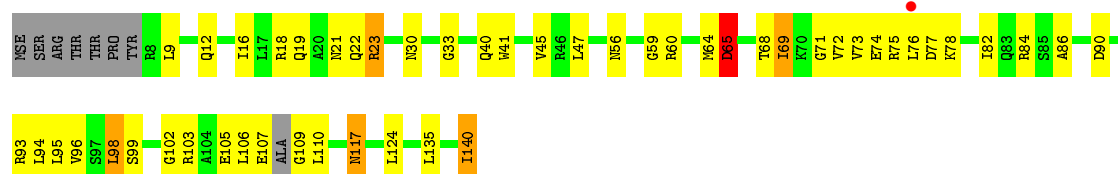
- Molecule 1: Probable transcriptional regulator

Chain F: 69% 22% 5%



- Molecule 1: Probable transcriptional regulator

Chain G: 59% 31% 6%



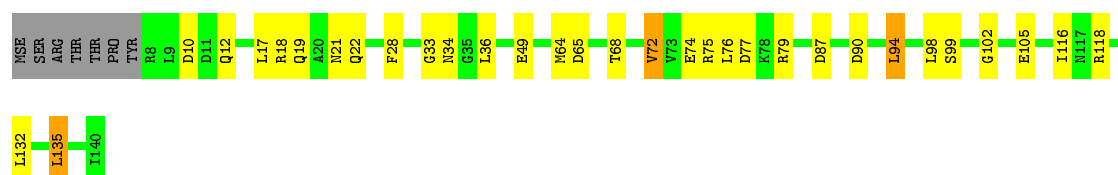
- Molecule 1: Probable transcriptional regulator

Chain H: 75% 19% 5%



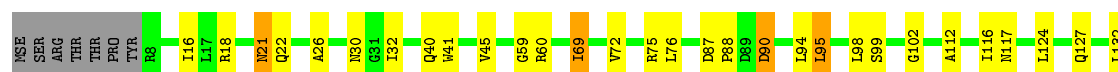
- Molecule 1: Probable transcriptional regulator

Chain I: 72% 21% 5%



- Molecule 1: Probable transcriptional regulator

Chain J: 71% 21% 5%







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	89.37Å 89.37Å 217.24Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.14 – 2.40 34.13 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.5 (34.14-2.40) 98.6 (34.13-2.40)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.09 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.218 , 0.262 0.217 , 0.260	Depositor DCC
$R_{free}$ test set	3761 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	36.9	Xtriage
Anisotropy	0.729	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 37.7	EDS
Estimated twinning fraction	0.010 for -h,-k,l 0.480 for h,-h-k,-l 0.014 for -k,-h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 74986 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	10506	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.59% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.49	0/1018	0.69	0/1374
1	B	0.49	0/1018	0.68	0/1374
1	C	0.44	0/1003	0.67	0/1352
1	D	0.48	0/1013	0.66	0/1367
1	E	0.43	0/1004	0.62	0/1353
1	F	0.45	0/1013	0.66	0/1367
1	G	0.46	0/1003	0.68	0/1352
1	H	0.48	0/1013	0.64	0/1367
1	I	0.45	0/1013	0.66	0/1367
1	J	0.43	0/1013	0.66	0/1367
All	All	0.46	0/10111	0.66	0/13640

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	1009	0	1035	34	0
1	B	1009	0	1035	37	0
1	C	995	0	1023	25	0
1	D	1004	0	1033	23	0
1	E	996	0	1028	44	0
1	F	1004	0	1033	42	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	995	0	1023	48	0
1	H	1004	0	1033	30	0
1	I	1004	0	1033	25	0
1	J	1004	0	1033	22	0
2	A	67	0	0	3	0
2	B	62	0	0	2	0
2	C	36	0	0	0	0
2	D	63	0	0	3	0
2	E	45	0	0	1	0
2	F	38	0	0	3	0
2	G	42	0	0	5	0
2	H	51	0	0	4	0
2	I	33	0	0	0	0
2	J	45	0	0	0	0
All	All	10506	0	10309	269	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 (269) 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:D:64:MSE:HE2	1:D:68:THR:HG22	1.19	1.15
1:B:64:MSE:HE2	1:B:68:THR:HG22	1.37	1.07
1:F:64:MSE:HE2	1:F:68:THR:HG22	1.29	1.06
1:C:64:MSE:HE2	1:C:68:THR:HG22	1.36	1.05
1:H:64:MSE:CE	1:H:68:THR:HG22	1.86	1.04
1:D:64:MSE:HE2	1:D:68:THR:CG2	1.88	1.03
1:B:64:MSE:HE1	1:B:72:VAL:CG1	1.88	1.03
1:B:64:MSE:HE1	1:B:72:VAL:HG11	1.36	1.03
1:H:64:MSE:HE1	1:H:72:VAL:CG1	1.90	1.01
1:F:64:MSE:CE	1:F:68:THR:HG22	1.91	1.00
1:I:64:MSE:HE2	1:I:68:THR:HG22	1.39	0.99
1:E:64:MSE:CE	1:E:68:THR:HG22	1.95	0.97
1:E:64:MSE:HE2	1:E:68:THR:HG22	1.41	0.97
1:B:64:MSE:CE	1:B:68:THR:HG22	1.95	0.94
1:G:64:MSE:HE2	1:G:68:THR:HG22	1.52	0.92
1:H:64:MSE:HE2	1:H:68:THR:HG22	1.52	0.91
1:H:64:MSE:HE3	1:H:68:THR:HG22	1.54	0.88
1:H:64:MSE:HE1	1:H:72:VAL:HG13	1.54	0.88
1:D:86:ALA:H	1:H:19:GLN:HE21	1.20	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:93:ARG:NH2	2:G:174:HOH:O	2.04	0.86
1:D:64:MSE:CE	1:D:68:THR:HG22	2.04	0.86
1:I:64:MSE:CE	1:I:68:THR:HG22	2.06	0.84
1:C:23:ARG:HH11	1:C:23:ARG:HG2	1.42	0.84
1:F:53:CYS:SG	2:F:168:HOH:O	2.33	0.83
1:B:86:ALA:H	1:I:19:GLN:HE21	1.26	0.83
1:H:64:MSE:HE1	1:H:72:VAL:HG11	1.62	0.81
1:E:140:ILE:HD11	1:F:16:ILE:HG23	1.62	0.81
1:G:64:MSE:CE	1:G:68:THR:HG22	2.11	0.80
1:A:19:GLN:HE21	1:F:86:ALA:H	1.24	0.80
1:E:40:GLN:HE22	1:E:75:ARG:HE	1.31	0.78
1:D:86:ALA:H	1:H:19:GLN:NE2	1.83	0.77
1:H:72:VAL:HA	1:H:75:ARG:HE	1.48	0.77
1:G:23:ARG:CG	1:G:23:ARG:HH11	2.00	0.75
1:E:40:GLN:NE2	1:E:75:ARG:HE	1.85	0.75
1:C:23:ARG:HH11	1:C:23:ARG:CG	2.02	0.73
1:C:64:MSE:HE2	1:C:68:THR:CG2	2.16	0.73
1:J:59:GLY:HA2	1:J:69:ILE:HG13	1.71	0.72
1:G:106:LEU:O	1:G:110:LEU:HG	1.91	0.71
1:F:64:MSE:HE1	1:F:72:VAL:CG1	2.20	0.71
1:B:140:ILE:HG22	1:F:94:LEU:HD11	1.71	0.71
1:C:137:ALA:HA	1:C:140:ILE:HD12	1.71	0.71
1:E:64:MSE:HE1	1:E:72:VAL:CG1	2.20	0.71
1:B:86:ALA:H	1:I:19:GLN:NE2	1.88	0.70
1:B:75:ARG:HD3	2:B:192:HOH:O	1.92	0.70
1:E:64:MSE:HE1	1:E:72:VAL:HG11	1.74	0.69
1:F:64:MSE:HE1	1:F:72:VAL:HG11	1.74	0.69
1:G:64:MSE:HE1	1:G:72:VAL:HG11	1.74	0.69
1:E:31:GLY:HA3	1:E:116:ILE:HD11	1.74	0.69
1:A:32:ILE:HD13	1:A:34:ASN:OD1	1.93	0.69
1:D:114:ARG:NH1	2:D:191:HOH:O	2.27	0.68
1:F:18:ARG:HH11	1:F:22:GLN:HE22	1.42	0.67
1:E:137:ALA:HA	1:E:140:ILE:HD13	1.78	0.66
1:C:77:ASP:HB2	1:C:82:ILE:HD11	1.78	0.65
1:G:77:ASP:HB2	1:G:82:ILE:HD11	1.79	0.65
1:J:87:ASP:HB3	1:J:90:ASP:O	1.97	0.64
1:B:72:VAL:HG12	1:B:75:ARG:HH21	1.62	0.64
1:G:33:GLY:HA2	1:I:33:GLY:HA2	1.79	0.64
1:H:90:ASP:OD1	1:H:92:ARG:HD3	1.98	0.64
1:J:40:GLN:HE21	1:J:75:ARG:HD2	1.62	0.64
1:F:140:ILE:HG22	1:G:94:LEU:CD1	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:40:GLN:NE2	1:J:75:ARG:HD2	2.13	0.63
1:C:139:LEU:HD21	1:D:124:LEU:HD22	1.81	0.62
1:I:99:SER:HB2	1:I:102:GLY:H	1.63	0.62
1:E:41:TRP:O	1:E:45:VAL:HG23	1.99	0.62
1:G:65:ASP:HB2	1:G:68:THR:OG1	1.99	0.62
1:I:19:GLN:HB2	1:J:140:ILE:HD12	1.80	0.62
1:G:16:ILE:HG23	1:H:140:ILE:HD11	1.82	0.62
1:D:97:SER:HB2	2:D:172:HOH:O	1.98	0.61
1:C:64:MSE:HE1	1:C:72:VAL:HG11	1.80	0.61
1:G:59:GLY:CA	1:G:69:ILE:HG13	2.30	0.61
1:E:21:ASN:HD22	1:F:17:LEU:HB3	1.66	0.61
1:C:64:MSE:CE	1:C:68:THR:HG22	2.23	0.61
1:E:59:GLY:HA2	1:E:69:ILE:HG13	1.83	0.61
1:F:99:SER:HB2	1:F:102:GLY:H	1.66	0.60
1:C:99:SER:HB2	1:C:102:GLY:H	1.67	0.60
1:D:18:ARG:HH11	1:D:22:GLN:HE22	1.47	0.60
1:F:18:ARG:HH11	1:F:22:GLN:NE2	2.00	0.60
1:C:124:LEU:HD13	1:D:139:LEU:HD21	1.83	0.59
1:H:18:ARG:HH11	1:H:22:GLN:HE22	1.51	0.59
1:G:124:LEU:HD13	1:H:139:LEU:HD21	1.83	0.59
1:B:64:MSE:HE3	1:B:68:THR:HG22	1.85	0.58
1:C:18:ARG:HH11	1:C:22:GLN:HE22	1.50	0.58
1:I:18:ARG:HH11	1:I:22:GLN:HE22	1.50	0.58
1:G:16:ILE:HG23	1:H:140:ILE:CD1	2.34	0.58
1:E:21:ASN:ND2	1:F:17:LEU:HB3	2.18	0.58
1:A:117:ASN:HD21	1:B:13:ILE:H	1.52	0.58
1:A:19:GLN:HE22	1:F:85:SER:HA	1.69	0.57
1:F:97:SER:HB3	2:F:178:HOH:O	2.03	0.57
1:G:23:ARG:HH11	1:G:23:ARG:HG2	1.69	0.57
1:G:23:ARG:HH11	1:G:23:ARG:HG3	1.70	0.57
1:E:64:MSE:HE2	1:E:68:THR:CG2	2.27	0.57
1:H:71:GLY:O	1:H:75:ARG:HD3	2.05	0.56
1:A:140:ILE:HG13	1:B:19:GLN:HB2	1.87	0.56
1:E:137:ALA:HA	1:E:140:ILE:CD1	2.36	0.56
1:I:64:MSE:HE2	1:I:68:THR:CG2	2.25	0.55
1:E:87:ASP:OD2	1:E:93:ARG:NE	2.36	0.55
1:B:60:ARG:NH1	1:F:52:PRO:HB2	2.22	0.55
1:F:64:MSE:HE2	1:F:68:THR:CG2	2.20	0.55
1:B:85:SER:HA	1:I:19:GLN:HE22	1.71	0.55
1:A:19:GLN:NE2	1:F:86:ALA:H	2.01	0.55
1:A:16:ILE:HG23	1:B:140:ILE:HD11	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:23:ARG:CG	1:G:23:ARG:NH1	2.66	0.54
1:B:72:VAL:HG12	1:B:75:ARG:NH2	2.23	0.54
1:A:97:SER:HB2	2:A:157:HOH:O	2.08	0.54
1:C:65:ASP:O	1:C:69:ILE:HG22	2.07	0.54
1:E:87:ASP:HB3	1:E:90:ASP:O	2.08	0.54
1:A:17:LEU:HB3	1:B:21:ASN:HD22	1.73	0.54
1:E:59:GLY:CA	1:E:69:ILE:HG13	2.38	0.54
1:E:64:MSE:HE3	1:E:68:THR:HG22	1.85	0.53
1:G:64:MSE:HE1	1:G:72:VAL:CG1	2.38	0.53
1:J:18:ARG:HH11	1:J:22:GLN:HE22	1.55	0.53
1:C:23:ARG:CG	1:C:23:ARG:NH1	2.70	0.53
1:F:18:ARG:NH1	1:F:22:GLN:HE22	2.04	0.53
1:E:64:MSE:HE1	1:E:72:VAL:HG13	1.91	0.53
1:E:87:ASP:HA	1:E:95:LEU:HD22	1.91	0.53
1:G:30:ASN:OD1	2:G:156:HOH:O	2.19	0.53
1:G:56:ASN:O	1:G:60:ARG:HG3	2.09	0.53
1:D:85:SER:HA	1:H:19:GLN:HE22	1.73	0.53
1:C:107:GLU:O	1:C:109:GLY:N	2.42	0.53
1:G:140:ILE:HD11	1:H:16:ILE:HA	1.91	0.53
1:F:115:GLU:O	1:F:119:GLN:HB3	2.09	0.52
1:H:12:GLN:O	1:H:16:ILE:HG12	2.09	0.52
1:D:72:VAL:HA	1:D:75:ARG:HE	1.73	0.52
1:E:121:LEU:HD21	1:F:13:ILE:HD12	1.92	0.52
1:I:64:MSE:HE1	1:I:72:VAL:CG1	2.40	0.52
1:C:72:VAL:O	1:C:76:LEU:HB2	2.09	0.52
1:J:18:ARG:HH11	1:J:22:GLN:NE2	2.08	0.52
1:I:28:PHE:HA	1:I:116:ILE:HD13	1.91	0.51
1:E:140:ILE:CD1	1:F:16:ILE:HG23	2.37	0.51
1:C:140:ILE:HD11	1:D:16:ILE:HG12	1.92	0.51
1:A:21:ASN:HD22	1:B:17:LEU:HB3	1.73	0.51
1:D:121:LEU:HD22	1:D:129:GLN:HG2	1.92	0.51
1:C:72:VAL:HG12	1:C:75:ARG:HH21	1.76	0.51
1:G:18:ARG:HH11	1:G:22:GLN:HE22	1.59	0.51
1:E:31:GLY:CA	1:E:116:ILE:HD11	2.41	0.51
1:G:23:ARG:HG2	1:G:23:ARG:NH1	2.26	0.51
1:E:139:LEU:HD21	1:F:124:LEU:HD22	1.92	0.51
1:B:114:ARG:NH1	2:B:201:HOH:O	2.44	0.51
1:G:40:GLN:HE21	1:G:75:ARG:HE	1.60	0.50
1:H:60:ARG:HD3	2:H:160:HOH:O	2.12	0.50
1:B:40:GLN:HE21	1:B:75:ARG:CZ	2.23	0.50
1:A:13:ILE:H	1:B:117:ASN:HD21	1.60	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:41:TRP:O	1:G:45:VAL:HG23	2.12	0.50
1:C:68:THR:O	1:C:72:VAL:HG13	2.12	0.50
1:F:57:GLN:HE22	1:G:103:ARG:HH22	1.60	0.49
1:H:97:SER:HB2	2:H:162:HOH:O	2.11	0.49
1:F:75:ARG:O	1:F:79:ARG:HG3	2.12	0.49
1:B:64:MSE:HE1	1:B:72:VAL:HG13	1.84	0.49
1:D:83:GLN:HG2	1:D:97:SER:HB3	1.93	0.49
1:E:60:ARG:HH11	1:E:60:ARG:HB3	1.78	0.49
1:D:52:PRO:O	1:G:60:ARG:NH1	2.44	0.49
1:B:18:ARG:HH11	1:B:22:GLN:HE22	1.59	0.49
1:J:26:ALA:O	1:J:30:ASN:ND2	2.43	0.49
1:G:73:VAL:HG12	1:G:84:ARG:HE	1.77	0.49
1:B:83:GLN:HG3	1:B:97:SER:HB3	1.94	0.49
1:E:19:GLN:HE21	1:G:86:ALA:H	1.60	0.49
1:A:18:ARG:HH11	1:A:22:GLN:NE2	2.11	0.49
1:B:18:ARG:HH11	1:B:22:GLN:NE2	2.10	0.48
1:A:140:ILE:HD11	1:B:16:ILE:HG23	1.94	0.48
1:B:115:GLU:OE1	1:B:118:ARG:HD3	2.14	0.48
1:E:31:GLY:HA3	1:E:116:ILE:CD1	2.43	0.48
1:I:18:ARG:NH1	1:I:22:GLN:HE22	2.09	0.48
1:I:64:MSE:HE1	1:I:72:VAL:HG11	1.96	0.48
1:E:69:ILE:HD13	1:E:73:VAL:HG23	1.94	0.48
1:C:18:ARG:HH11	1:C:22:GLN:NE2	2.11	0.48
1:I:28:PHE:HE1	1:I:36:LEU:O	1.96	0.48
1:G:64:MSE:HE2	1:G:68:THR:CG2	2.34	0.48
1:E:19:GLN:NE2	1:G:86:ALA:H	2.12	0.48
1:J:112:ALA:O	1:J:116:ILE:HG12	2.13	0.48
1:A:115:GLU:HA	1:A:118:ARG:HG2	1.95	0.47
1:A:117:ASN:ND2	1:B:13:ILE:H	2.12	0.47
1:H:40:GLN:HE21	1:H:75:ARG:CZ	2.27	0.47
1:J:87:ASP:HA	1:J:95:LEU:HD22	1.95	0.47
1:C:56:ASN:O	1:C:60:ARG:HG3	2.15	0.47
1:J:132:LEU:O	1:J:136:LEU:HB2	2.14	0.47
1:G:84:ARG:NH1	2:G:164:HOH:O	2.44	0.47
1:C:106:LEU:O	1:C:110:LEU:HG	2.14	0.47
1:F:64:MSE:HE3	1:F:68:THR:HG22	1.89	0.47
1:A:28:PHE:CZ	1:A:32:ILE:HD12	2.50	0.47
1:G:117:ASN:HB2	1:H:13:ILE:HG22	1.95	0.47
1:D:53:CYS:SG	1:D:57:GLN:HG2	2.55	0.47
1:I:18:ARG:HH11	1:I:22:GLN:NE2	2.12	0.47
1:G:107:GLU:O	1:G:109:GLY:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:23:ARG:HG2	1:C:23:ARG:NH1	2.19	0.46
1:H:60:ARG:CD	2:H:160:HOH:O	2.63	0.46
1:G:71:GLY:O	1:G:75:ARG:HD3	2.15	0.46
1:A:93:ARG:HB2	2:A:194:HOH:O	2.16	0.46
1:C:82:ILE:HD13	1:C:96:VAL:HG13	1.98	0.46
1:A:140:ILE:HG13	1:B:19:GLN:CB	2.46	0.46
1:A:18:ARG:HH11	1:A:22:GLN:HE22	1.62	0.46
1:A:28:PHE:HA	1:A:116:ILE:HD13	1.98	0.46
1:J:140:ILE:HA	1:J:140:ILE:HD13	1.66	0.45
1:G:19:GLN:CB	1:H:140:ILE:HG12	2.46	0.45
1:A:140:ILE:CD1	1:B:16:ILE:HG23	2.46	0.45
1:H:8:ARG:N	2:H:165:HOH:O	2.48	0.45
1:I:12:GLN:OE1	1:J:117:ASN:ND2	2.49	0.45
1:A:47:LEU:CB	1:A:98:LEU:HD13	2.47	0.45
1:F:72:VAL:O	1:F:76:LEU:HB2	2.17	0.45
1:E:40:GLN:NE2	1:E:75:ARG:NE	2.60	0.45
1:F:140:ILE:HG22	1:G:94:LEU:HD11	1.97	0.45
1:H:64:MSE:HE3	1:H:68:THR:CG2	2.36	0.45
1:I:135:LEU:HB3	1:J:132:LEU:HD13	1.99	0.45
1:E:17:LEU:HB3	1:F:21:ASN:HD22	1.81	0.45
1:D:64:MSE:HE1	1:D:72:VAL:HG11	1.99	0.44
1:E:121:LEU:CD2	1:F:13:ILE:HD12	2.47	0.44
1:G:117:ASN:C	1:G:117:ASN:OD1	2.56	0.44
1:E:132:LEU:O	1:E:136:LEU:HB2	2.17	0.44
1:J:41:TRP:O	1:J:45:VAL:HG23	2.17	0.44
1:D:65:ASP:HB2	2:D:150:HOH:O	2.17	0.44
1:A:87:ASP:HB3	1:A:93:ARG:HB3	1.99	0.44
1:I:87:ASP:HB3	1:I:90:ASP:O	2.17	0.44
1:F:76:LEU:HB3	1:F:82:ILE:HG12	2.00	0.44
1:B:140:ILE:HG22	1:F:94:LEU:CD1	2.46	0.44
1:G:30:ASN:ND2	2:G:146:HOH:O	2.50	0.44
1:H:64:MSE:HE2	1:H:68:THR:C	2.38	0.44
1:B:52:PRO:O	1:J:60:ARG:NH1	2.51	0.43
1:A:66:ALA:O	1:A:70:LYS:HB2	2.18	0.43
1:G:9:LEU:HD12	1:G:12:GLN:HG3	2.00	0.43
1:G:93:ARG:CZ	2:G:174:HOH:O	2.60	0.43
1:I:17:LEU:HB3	1:J:21:ASN:ND2	2.34	0.43
1:E:47:LEU:HB3	1:E:98:LEU:HD13	2.00	0.43
1:I:132:LEU:HD21	1:J:139:LEU:HD11	1.99	0.43
1:G:56:ASN:HB2	1:G:93:ARG:HH22	1.84	0.43
1:A:115:GLU:OE1	1:A:118:ARG:HD3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:65:ASP:HB2	1:I:68:THR:H	1.83	0.43
1:G:82:ILE:HD13	1:G:96:VAL:HG13	2.01	0.43
1:A:65:ASP:HB3	1:A:68:THR:H	1.84	0.43
1:A:13:ILE:H	1:B:117:ASN:ND2	2.15	0.43
1:A:133:ARG:NH2	2:A:178:HOH:O	2.52	0.43
1:A:16:ILE:HG23	1:B:140:ILE:CD1	2.49	0.42
1:E:135:LEU:HB3	1:F:132:LEU:HD13	2.01	0.42
1:J:59:GLY:CA	1:J:69:ILE:HG13	2.46	0.42
1:E:27:LEU:O	1:E:116:ILE:HD12	2.19	0.42
1:E:40:GLN:HE22	1:E:75:ARG:NE	2.06	0.42
1:I:94:LEU:HD12	1:I:94:LEU:HA	1.79	0.42
1:F:30:ASN:HB2	2:F:169:HOH:O	2.19	0.42
1:A:47:LEU:HB3	1:A:98:LEU:HD13	2.02	0.42
1:F:140:ILE:CG2	1:G:94:LEU:HD11	2.49	0.42
1:I:34:ASN:HB2	1:I:105:GLU:OE2	2.20	0.42
1:F:87:ASP:HB3	1:F:90:ASP:O	2.19	0.42
1:G:140:ILE:HD12	1:G:140:ILE:HA	1.77	0.42
1:J:87:ASP:HA	1:J:95:LEU:CD2	2.50	0.42
1:H:57:GLN:OE1	1:H:60:ARG:NH1	2.53	0.42
1:H:115:GLU:OE1	1:H:118:ARG:NH1	2.50	0.42
1:F:115:GLU:OE1	1:F:118:ARG:NH1	2.53	0.42
1:E:18:ARG:HH11	1:E:22:GLN:HE22	1.67	0.42
1:J:99:SER:HB2	1:J:102:GLY:H	1.85	0.41
1:C:121:LEU:HD22	1:C:129:GLN:HG2	2.02	0.41
1:G:99:SER:HB2	1:G:102:GLY:H	1.86	0.41
1:A:40:GLN:HE21	1:A:75:ARG:CZ	2.34	0.41
1:E:112:ALA:O	1:E:116:ILE:HG12	2.21	0.41
1:D:94:LEU:HD11	1:G:140:ILE:HG22	2.03	0.41
1:I:77:ASP:C	1:I:79:ARG:H	2.24	0.41
1:D:64:MSE:HE2	1:D:68:THR:HG21	1.93	0.41
1:A:19:GLN:NE2	1:F:85:SER:HA	2.34	0.41
1:J:87:ASP:HA	1:J:88:PRO:HD2	1.92	0.41
1:B:60:ARG:HH12	1:F:52:PRO:HB2	1.86	0.41
1:F:87:ASP:HA	1:F:95:LEU:HD22	2.03	0.41
1:A:132:LEU:HD13	1:B:135:LEU:HB3	2.03	0.41
1:F:47:LEU:HB2	1:F:98:LEU:HD13	2.02	0.41
1:E:64:MSE:CE	1:E:68:THR:CG2	2.84	0.41
1:D:116:ILE:HA	1:D:119:GLN:HE21	1.86	0.41
1:E:106:LEU:O	1:E:110:LEU:HG	2.22	0.40
1:A:19:GLN:HB2	1:B:140:ILE:HG13	2.02	0.40
1:E:18:ARG:HH11	1:E:22:GLN:NE2	2.19	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:122:ALA:N	1:H:123:PRO:CD	2.85	0.40
1:D:41:TRP:O	1:D:45:VAL:HG23	2.21	0.40
1:E:64:MSE:SE	2:E:185:HOH:O	2.90	0.40
1:G:47:LEU:CB	1:G:98:LEU:HD13	2.52	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	132/140 (94%)	129 (98%)	3 (2%)	0	100	100
1	B	132/140 (94%)	130 (98%)	2 (2%)	0	100	100
1	C	128/140 (91%)	125 (98%)	3 (2%)	0	100	100
1	D	131/140 (94%)	131 (100%)	0	0	100	100
1	E	128/140 (91%)	127 (99%)	1 (1%)	0	100	100
1	F	131/140 (94%)	131 (100%)	0	0	100	100
1	G	128/140 (91%)	126 (98%)	0	2 (2%)	12	16
1	H	131/140 (94%)	130 (99%)	1 (1%)	0	100	100
1	I	131/140 (94%)	128 (98%)	3 (2%)	0	100	100
1	J	131/140 (94%)	130 (99%)	0	1 (1%)	24	35
All	All	1303/1400 (93%)	1287 (99%)	13 (1%)	3 (0%)	52	69

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	J	90	ASP
1	G	65	ASP
1	G	90	ASP

### 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	100/106 (94%)	90 (90%)	10 (10%)	9	14
1	B	100/106 (94%)	90 (90%)	10 (10%)	9	14
1	C	99/106 (93%)	89 (90%)	10 (10%)	9	13
1	D	100/106 (94%)	91 (91%)	9 (9%)	12	17
1	E	99/106 (93%)	86 (87%)	13 (13%)	5	6
1	F	100/106 (94%)	91 (91%)	9 (9%)	12	17
1	G	99/106 (93%)	86 (87%)	13 (13%)	5	6
1	H	100/106 (94%)	94 (94%)	6 (6%)	24	37
1	I	100/106 (94%)	89 (89%)	11 (11%)	8	10
1	J	100/106 (94%)	87 (87%)	13 (13%)	5	6
All	All	997/1060 (94%)	893 (90%)	104 (10%)	9	12

All (104) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	32	ILE
1	A	65	ASP
1	A	76	LEU
1	A	94	LEU
1	A	95	LEU
1	A	98	LEU
1	A	124	LEU
1	A	135	LEU
1	A	140	ILE
1	B	21	ASN
1	B	69	ILE
1	B	72	VAL
1	B	76	LEU
1	B	94	LEU
1	B	95	LEU

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Mol	Chain	Res	Type
1	B	98	LEU
1	B	124	LEU
1	B	135	LEU
1	B	140	ILE
1	C	10	ASP
1	C	23	ARG
1	C	72	VAL
1	C	76	LEU
1	C	78	LYS
1	C	83	GLN
1	C	95	LEU
1	C	98	LEU
1	C	135	LEU
1	C	140	ILE
1	D	57	GLN
1	D	75	ARG
1	D	76	LEU
1	D	94	LEU
1	D	95	LEU
1	D	98	LEU
1	D	124	LEU
1	D	127	GLN
1	D	138	ARG
1	E	21	ASN
1	E	60	ARG
1	E	69	ILE
1	E	76	LEU
1	E	83	GLN
1	E	95	LEU
1	E	98	LEU
1	E	118	ARG
1	E	124	LEU
1	E	127	GLN
1	E	135	LEU
1	E	136	LEU
1	E	140	ILE
1	F	10	ASP
1	F	13	ILE
1	F	21	ASN
1	F	72	VAL
1	F	76	LEU
1	F	95	LEU

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Mol	Chain	Res	Type
1	F	103	ARG
1	F	124	LEU
1	F	135	LEU
1	G	21	ASN
1	G	23	ARG
1	G	65	ASP
1	G	69	ILE
1	G	74	GLU
1	G	76	LEU
1	G	78	LYS
1	G	95	LEU
1	G	98	LEU
1	G	105	GLU
1	G	117	ASN
1	G	135	LEU
1	G	140	ILE
1	H	72	VAL
1	H	75	ARG
1	H	76	LEU
1	H	94	LEU
1	H	95	LEU
1	H	98	LEU
1	I	10	ASP
1	I	21	ASN
1	I	49	GLU
1	I	72	VAL
1	I	74	GLU
1	I	75	ARG
1	I	76	LEU
1	I	94	LEU
1	I	98	LEU
1	I	118	ARG
1	I	135	LEU
1	J	16	ILE
1	J	21	ASN
1	J	32	ILE
1	J	69	ILE
1	J	72	VAL
1	J	76	LEU
1	J	94	LEU
1	J	95	LEU
1	J	98	LEU

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Mol	Chain	Res	Type
1	J	124	LEU
1	J	127	GLN
1	J	135	LEU
1	J	136	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (50) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	21	ASN
1	A	22	GLN
1	A	40	GLN
1	A	55	GLN
1	A	83	GLN
1	A	117	ASN
1	A	119	GLN
1	B	21	ASN
1	B	22	GLN
1	B	40	GLN
1	B	117	ASN
1	B	119	GLN
1	C	12	GLN
1	C	21	ASN
1	C	22	GLN
1	C	83	GLN
1	D	22	GLN
1	D	55	GLN
1	D	83	GLN
1	D	117	ASN
1	D	119	GLN
1	E	19	GLN
1	E	21	ASN
1	E	22	GLN
1	E	40	GLN
1	E	83	GLN
1	E	117	ASN
1	E	119	GLN
1	F	21	ASN
1	F	22	GLN
1	G	21	ASN
1	G	22	GLN
1	H	19	GLN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	H	22	GLN
1	H	40	GLN
1	H	55	GLN
1	H	83	GLN
1	H	119	GLN
1	I	19	GLN
1	I	21	ASN
1	I	22	GLN
1	I	55	GLN
1	I	83	GLN
1	J	12	GLN
1	J	21	ASN
1	J	22	GLN
1	J	83	GLN
1	J	117	ASN
1	J	119	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 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

There are no chain breaks in this entry.



## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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(Å²)	Q<0.9	
1	A	133/140 (95%)	-0.11	0	100	100	28, 42, 52, 57	0
1	B	133/140 (95%)	-0.08	0	100	100	28, 42, 52, 57	0
1	C	131/140 (93%)	0.04	0	100	100	37, 49, 62, 63	0
1	D	132/140 (94%)	-0.08	0	100	100	28, 42, 58, 62	0
1	E	131/140 (93%)	-0.00	0	100	100	42, 53, 62, 64	0
1	F	132/140 (94%)	0.08	0	100	100	35, 46, 60, 64	0
1	G	131/140 (93%)	0.08	1 (0%)	87	87	37, 49, 61, 65	0
1	H	132/140 (94%)	-0.15	0	100	100	31, 42, 60, 65	0
1	I	132/140 (94%)	0.02	0	100	100	35, 47, 61, 65	0
1	J	132/140 (94%)	0.03	0	100	100	41, 54, 62, 65	0
All	All	1319/1400 (94%)	-0.02	1 (0%)	95	95	28, 47, 61, 65	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	76	LEU	2.2

### 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.