



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 01:00 am BST

PDB ID : 3DKU
Title : Crystal structure of Nudix hydrolase Orf153, ymfB, from Escherichia coli K-1
Authors : Hong, M.K.; Kim, J.K.; Jung, J.H.; Jung, J.W.; Choi, J.Y.; Kang, L.W.
Deposited on : 2008-06-26
Resolution : 2.69 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
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

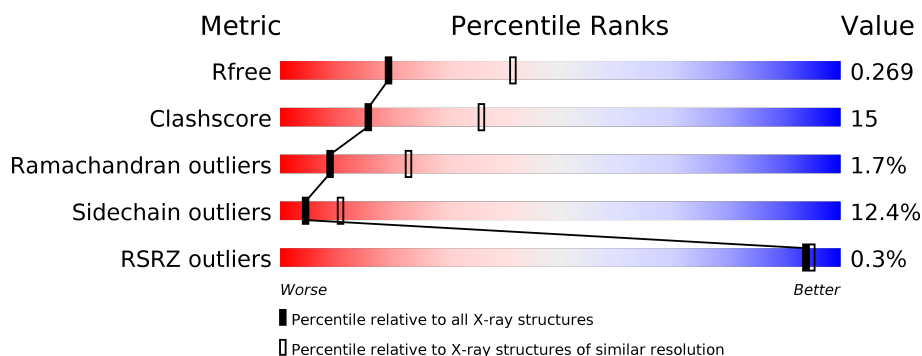
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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 ≥ 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	153	<div> <div>%</div> <div> <div></div> <div>63%</div> <div>31%</div> <div>5%</div> </div> </div>
1	B	153	<div> <div></div> <div>58%</div> <div>32%</div> <div>5%</div> <div>• •</div> </div>
1	C	153	<div> <div>%</div> <div> <div></div> <div>70%</div> <div>25%</div> <div>5%</div> </div> </div>
1	D	153	<div> <div></div> <div>67%</div> <div>24%</div> <div>5%</div> <div>• •</div> </div>
1	E	153	<div> <div></div> <div>62%</div> <div>30%</div> <div>7%</div> <div>•</div> </div>
1	F	153	<div> <div></div> <div>65%</div> <div>27%</div> <div>5%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	153	<div><div></div><div>64%27%8%</div><div></div></div>
1	H	153	<div>%<div><div></div><div>59%29%6%</div><div></div></div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9800 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 Putative phosphohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	0	0	0
			1227	786	209	224	8			
1	B	147	Total	C	N	O	S	0	0	0
			1181	754	202	217	8			
1	C	153	Total	C	N	O	S	0	0	0
			1227	786	209	224	8			
1	D	148	Total	C	N	O	S	0	0	0
			1192	763	203	218	8			
1	E	152	Total	C	N	O	S	0	0	0
			1219	780	208	223	8			
1	F	149	Total	C	N	O	S	0	0	0
			1199	767	204	220	8			
1	G	152	Total	C	N	O	S	0	0	0
			1219	780	208	223	8			
1	H	147	Total	C	N	O	S	0	0	0
			1181	754	202	217	8			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	18	Total	O	0	0
			18	18		
2	B	19	Total	O	0	0
			19	19		
2	C	27	Total	O	0	0
			27	27		
2	D	20	Total	O	0	0
			20	20		
2	E	20	Total	O	0	0
			20	20		
2	F	22	Total	O	0	0
			22	22		

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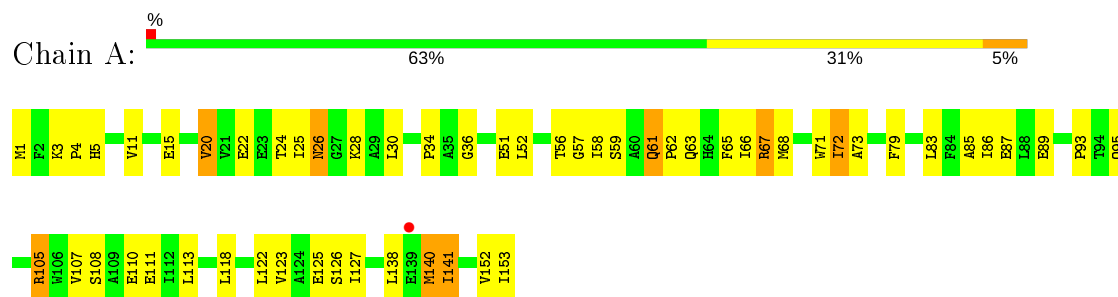
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	9	Total	O	0	0
			9	9		
2	H	20	Total	O	0	0
			20	20		

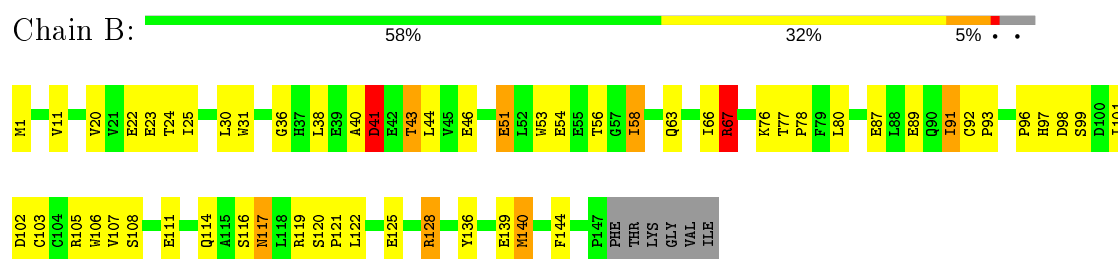
3 Residue-property plots

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

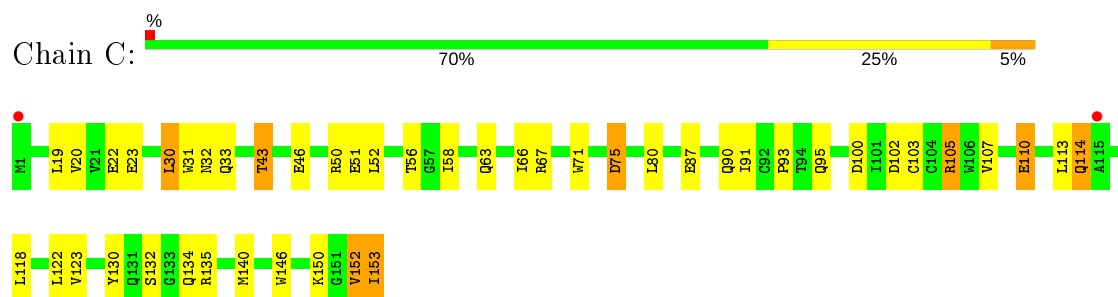
• Molecule 1: Putative phosphohydrolase



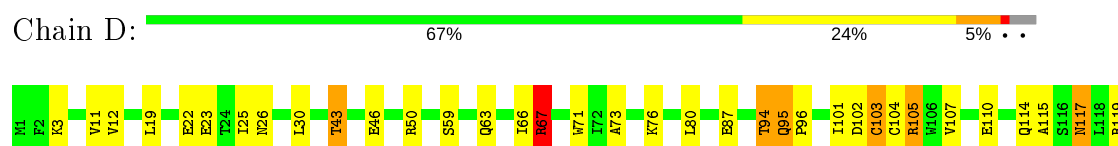
• Molecule 1: Putative phosphohydrolase

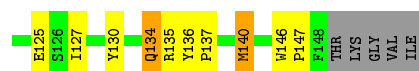


• Molecule 1: Putative phosphohydrolase



• Molecule 1: Putative phosphohydrolase





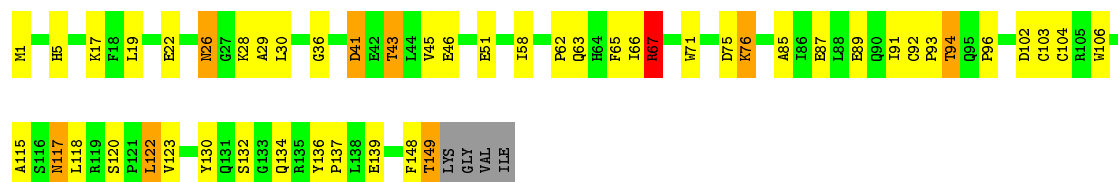
- Molecule 1: Putative phosphohydrolase

Chain E: 62% 30% 7% •



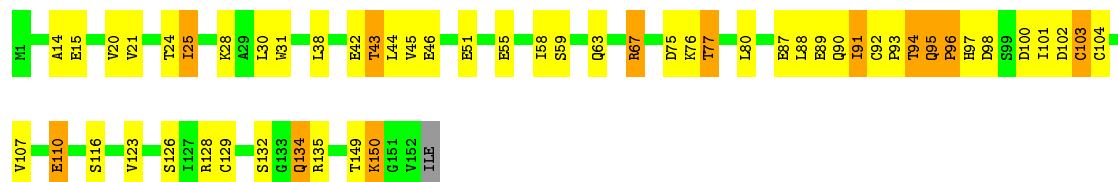
- Molecule 1: Putative phosphohydrolase

Chain F: 65% 27% 5% • •



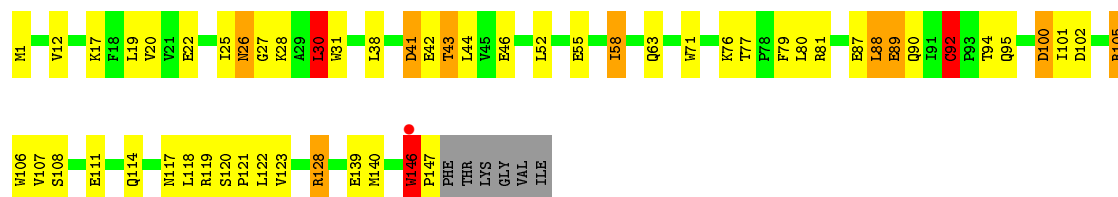
- Molecule 1: Putative phosphohydrolase

Chain G: 64% 27% 8% •



- Molecule 1: Putative phosphohydrolase

Chain H: 59% 29% 6% • •



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.96Å 70.52Å 145.03Å 90.00° 103.30° 90.00°	Depositor
Resolution (Å)	33.56 – 2.69 33.56 – 2.69	Depositor EDS
% Data completeness (in resolution range)	98.3 (33.56-2.69) 98.3 (33.56-2.69)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.60 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.193 , 0.270 0.194 , 0.269	Depositor DCC
R_{free} test set	1865 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	33.9	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.019 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9800	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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	0/1262	0.84	0/1719
1	B	0.90	2/1215 (0.2%)	0.89	1/1656 (0.1%)
1	C	0.83	0/1262	0.85	2/1719 (0.1%)
1	D	0.86	0/1227	0.93	3/1672 (0.2%)
1	E	0.92	1/1254 (0.1%)	0.84	1/1708 (0.1%)
1	F	0.94	2/1234 (0.2%)	0.83	0/1682
1	G	0.89	1/1254 (0.1%)	0.88	2/1708 (0.1%)
1	H	0.92	1/1215 (0.1%)	0.92	4/1656 (0.2%)
All	All	0.89	7/9923 (0.1%)	0.87	13/13520 (0.1%)

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	H	0	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	92	CYS	CB-SG	-11.05	1.63	1.82
1	G	129	CYS	CB-SG	-6.64	1.71	1.82
1	E	139	GLU	CG-CD	6.52	1.61	1.51
1	B	139	GLU	CG-CD	5.61	1.60	1.51
1	B	51	GLU	CG-CD	5.52	1.60	1.51
1	F	139	GLU	CG-CD	5.29	1.59	1.51
1	F	104	CYS	CB-SG	-5.05	1.73	1.81

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	146	TRP	C-N-CD	-7.83	103.37	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	146	TRP	C-N-CD	-7.48	104.15	120.60
1	H	146	TRP	C-N-CA	6.32	148.54	122.00
1	H	30	LEU	CA-CB-CG	6.19	129.53	115.30
1	G	67	ARG	NE-CZ-NH1	-6.06	117.27	120.30
1	D	50	ARG	NE-CZ-NH2	-5.95	117.33	120.30
1	D	146	TRP	C-N-CA	5.68	145.87	122.00
1	H	88	LEU	CA-CB-CG	5.65	128.29	115.30
1	B	128	ARG	NE-CZ-NH2	-5.23	117.69	120.30
1	G	30	LEU	CA-CB-CG	5.19	127.24	115.30
1	C	75	ASP	N-CA-CB	-5.18	101.28	110.60
1	E	140	MET	CB-CG-SD	-5.11	97.06	112.40
1	C	30	LEU	CA-CB-CG	5.03	126.87	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	146	TRP	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	1227	0	1190	45	0
1	B	1181	0	1138	38	0
1	C	1227	0	1190	22	0
1	D	1192	0	1147	30	0
1	E	1219	0	1179	43	0
1	F	1199	0	1154	38	0
1	G	1219	0	1179	36	0
1	H	1181	0	1138	42	0
2	A	18	0	0	0	0
2	B	19	0	0	2	0
2	C	27	0	0	1	0
2	D	20	0	0	0	0
2	E	20	0	0	3	0
2	F	22	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	9	0	0	0	0
2	H	20	0	0	2	0
All	All	9800	0	9315	278	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 15.

All (278) 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:25:ILE:HG23	1:A:26:ASN:H	1.11	1.06
1:B:36:GLY:HA3	1:B:51:GLU:HG3	1.38	1.06
1:G:58:ILE:HD12	1:G:92:CYS:SG	2.02	0.99
1:H:43:THR:HG22	1:H:46:GLU:H	1.25	0.98
1:F:63:GLN:NE2	1:F:87:GLU:H	1.65	0.95
1:G:43:THR:HG22	1:G:46:GLU:H	1.33	0.94
1:F:43:THR:HG22	1:F:46:GLU:H	1.30	0.93
1:A:25:ILE:CG2	1:A:26:ASN:H	1.82	0.92
1:F:63:GLN:HE22	1:F:87:GLU:H	0.96	0.92
1:B:63:GLN:NE2	1:B:87:GLU:H	1.68	0.92
1:A:22:GLU:OE2	1:A:105:ARG:NH1	2.02	0.92
1:F:41:ASP:HB2	1:F:148:PHE:CE2	2.06	0.90
1:A:25:ILE:HG23	1:A:26:ASN:N	1.86	0.90
1:G:58:ILE:CD1	1:G:92:CYS:SG	2.63	0.86
1:B:36:GLY:HA3	1:B:51:GLU:CG	2.08	0.83
1:C:31:TRP:HE1	1:C:105:ARG:HH21	1.21	0.83
1:E:21:VAL:HG12	1:E:104:CYS:HB3	1.58	0.83
1:D:43:THR:HG22	1:D:46:GLU:H	1.45	0.81
1:G:110:GLU:CD	1:G:110:GLU:H	1.84	0.80
1:H:107:VAL:HG23	1:H:111:GLU:OE1	1.82	0.79
1:E:43:THR:HG22	1:E:46:GLU:H	1.47	0.79
1:E:20:VAL:HG13	1:E:31:TRP:CE3	2.21	0.76
1:D:22:GLU:OE2	1:D:105:ARG:NH1	2.18	0.76
1:D:66:ILE:O	1:D:67:ARG:HB2	1.85	0.75
1:G:102:ASP:O	1:G:103:CYS:HB3	1.85	0.74
1:F:63:GLN:HE22	1:F:87:GLU:N	1.81	0.73
1:F:66:ILE:O	1:F:67:ARG:HB2	1.89	0.73
1:G:63:GLN:HE22	1:G:87:GLU:H	1.37	0.73
1:B:63:GLN:HE22	1:B:87:GLU:H	1.36	0.73
1:C:110:GLU:H	1:C:110:GLU:CD	1.92	0.72
1:C:22:GLU:OE2	1:C:105:ARG:NH2	2.23	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:PHE:CD2	1:A:140:MET:HE1	2.24	0.71
1:H:139:GLU:HG3	1:H:140:MET:CE	2.20	0.71
1:H:107:VAL:HG22	1:H:108:SER:N	2.06	0.71
1:G:132:SER:OG	1:G:134:GLN:HG3	1.90	0.70
1:B:117:ASN:HD22	1:B:117:ASN:H	1.40	0.70
1:H:146:TRP:HA	1:H:146:TRP:CE3	2.27	0.70
1:B:136:TYR:HB3	1:B:140:MET:HE3	1.71	0.70
1:F:41:ASP:HB2	1:F:148:PHE:HE2	1.53	0.70
1:H:58:ILE:HD12	1:H:88:LEU:HD11	1.73	0.69
1:D:117:ASN:HD22	1:D:117:ASN:N	1.90	0.69
1:G:51:GLU:O	1:G:55:GLU:HG3	1.94	0.68
1:B:96:PRO:HB3	1:B:101:ILE:HB	1.75	0.68
1:H:146:TRP:HE3	1:H:146:TRP:HA	1.60	0.67
1:E:58:ILE:HD12	1:E:92:CYS:SG	2.34	0.67
1:F:75:ASP:O	1:F:76:LYS:HG2	1.94	0.66
1:D:117:ASN:HD22	1:D:117:ASN:H	1.43	0.66
1:A:66:ILE:HD11	1:A:85:ALA:HB2	1.78	0.66
1:H:100:ASP:N	1:H:100:ASP:OD2	2.25	0.65
1:B:128:ARG:HD2	2:B:173:HOH:O	1.95	0.65
1:C:63:GLN:NE2	1:C:87:GLU:H	1.95	0.65
1:A:72:ILE:HD12	1:A:73:ALA:O	1.97	0.65
1:A:67:ARG:HG2	1:A:68:MET:H	1.60	0.65
1:G:25:ILE:O	1:G:25:ILE:HG13	1.96	0.65
1:B:43:THR:HG22	1:B:46:GLU:HB2	1.79	0.64
1:G:46:GLU:OE1	1:G:150:LYS:HB3	1.97	0.64
1:C:146:TRP:CD1	1:C:150:LYS:HD2	2.33	0.63
1:D:43:THR:HB	1:D:46:GLU:OE2	1.98	0.62
1:H:107:VAL:CG2	1:H:108:SER:N	2.63	0.62
1:B:92:CYS:HB2	1:B:93:PRO:CD	2.30	0.62
1:E:75:ASP:O	1:E:76:LYS:CB	2.47	0.62
1:H:43:THR:CG2	1:H:46:GLU:H	2.06	0.61
1:E:31:TRP:CH2	1:E:105:ARG:HG2	2.35	0.61
1:H:139:GLU:HG3	1:H:140:MET:HE3	1.82	0.61
1:A:66:ILE:O	1:A:67:ARG:HB2	1.99	0.61
1:D:94:THR:HG21	1:D:104:CYS:HB2	1.82	0.61
1:G:46:GLU:OE1	1:G:150:LYS:CB	2.48	0.60
1:E:93:PRO:HA	1:G:95:GLN:HB2	1.83	0.60
1:H:63:GLN:NE2	1:H:87:GLU:H	1.99	0.60
1:F:117:ASN:HD22	1:F:117:ASN:H	1.48	0.60
1:A:67:ARG:HG2	1:A:68:MET:N	2.16	0.60
1:E:94:THR:HB	2:E:167:HOH:O	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:GLN:HE22	1:C:87:GLU:H	1.49	0.59
1:F:117:ASN:ND2	1:F:117:ASN:H	2.00	0.59
1:F:36:GLY:HA3	1:F:51:GLU:HG3	1.83	0.59
1:A:20:VAL:HG12	1:A:105:ARG:HG2	1.83	0.59
1:E:19:LEU:HD21	1:E:56:THR:HG21	1.84	0.59
1:A:107:VAL:HG22	1:A:111:GLU:HB2	1.83	0.59
1:A:125:GLU:HG3	1:B:144:PHE:CE1	2.38	0.59
1:A:93:PRO:HA	1:C:95:GLN:HB2	1.84	0.59
1:H:43:THR:HG22	1:H:46:GLU:N	2.08	0.59
1:H:26:ASN:O	1:H:28:LYS:N	2.33	0.58
1:A:63:GLN:NE2	1:A:87:GLU:H	2.01	0.58
1:F:17:LYS:HD3	1:F:91:ILE:HD11	1.84	0.58
1:H:139:GLU:HG3	1:H:140:MET:HE2	1.85	0.58
1:G:128:ARG:HH21	1:G:128:ARG:HG2	1.69	0.58
1:G:63:GLN:NE2	1:G:87:GLU:H	2.02	0.58
1:H:30:LEU:HA	1:H:117:ASN:O	2.04	0.57
1:D:94:THR:CG2	1:D:104:CYS:HB2	2.34	0.57
1:E:75:ASP:O	1:E:76:LYS:HB2	2.04	0.57
1:C:19:LEU:HD21	1:C:56:THR:HG21	1.86	0.57
1:B:53:TRP:CZ2	1:B:97:HIS:ND1	2.72	0.56
1:D:30:LEU:HD23	1:D:119:ARG:HA	1.87	0.56
1:B:66:ILE:O	1:B:67:ARG:HB2	2.02	0.56
1:F:28:LYS:HA	1:F:28:LYS:HE2	1.87	0.56
1:D:43:THR:HG22	1:D:46:GLU:HB2	1.87	0.56
1:E:25:ILE:HD12	1:E:119:ARG:HH21	1.70	0.56
1:E:105:ARG:HB3	1:G:103:CYS:HB3	1.88	0.56
1:E:20:VAL:HG13	1:E:31:TRP:HE3	1.70	0.56
1:A:125:GLU:HG3	1:B:144:PHE:CZ	2.41	0.56
1:C:102:ASP:O	1:C:103:CYS:HB3	2.05	0.56
1:G:58:ILE:HD11	1:G:92:CYS:SG	2.46	0.56
1:H:25:ILE:HG22	1:H:30:LEU:CD2	2.36	0.56
1:G:92:CYS:HB2	1:G:93:PRO:HD2	1.89	0.55
1:B:36:GLY:HA3	1:B:51:GLU:CD	2.27	0.55
1:A:95:GLN:HB2	1:C:93:PRO:HA	1.89	0.54
1:H:128:ARG:NE	2:H:166:HOH:O	2.41	0.54
1:H:63:GLN:NE2	1:H:87:GLU:N	2.55	0.54
1:A:56:THR:C	1:A:58:ILE:H	2.11	0.54
1:D:43:THR:CG2	1:D:46:GLU:H	2.18	0.54
1:D:66:ILE:O	1:D:67:ARG:CB	2.53	0.53
1:E:73:ALA:HB1	1:E:74:PRO:HD2	1.90	0.53
1:E:38:LEU:HD13	1:E:44:LEU:HD23	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:3:LYS:HD3	1:F:5:HIS:CD2	2.43	0.53
1:F:19:LEU:HD13	1:F:106:TRP:CE2	2.43	0.53
1:E:76:LYS:HE3	1:E:76:LYS:HA	1.89	0.53
1:H:118:LEU:HD22	1:H:123:VAL:HG11	1.89	0.53
1:A:26:ASN:HD22	1:A:26:ASN:N	2.06	0.53
1:B:63:GLN:HE22	1:B:87:GLU:N	2.06	0.53
1:G:123:VAL:O	1:G:126:SER:HB2	2.09	0.53
1:G:75:ASP:OD2	1:G:77:THR:OG1	2.28	0.52
1:B:22:GLU:OE2	1:B:31:TRP:NE1	2.42	0.52
1:E:1:MET:O	1:E:3:LYS:HE2	2.09	0.52
1:H:38:LEU:HD13	1:H:44:LEU:HD23	1.90	0.52
1:G:128:ARG:HH21	1:G:128:ARG:CG	2.23	0.52
1:F:102:ASP:O	1:F:103:CYS:HB3	2.09	0.52
1:C:71:TRP:CD1	1:C:122:LEU:HD11	2.46	0.51
1:D:135:ARG:C	1:D:136:TYR:CD2	2.84	0.51
1:A:62:PRO:HA	1:A:86:ILE:HG12	1.92	0.51
1:B:23:GLU:HG2	1:B:24:THR:N	2.26	0.51
1:C:32:ASN:OD1	1:C:33:GLN:N	2.43	0.51
1:G:96:PRO:O	1:G:97:HIS:HB2	2.10	0.51
1:E:131:GLN:HB3	1:H:90:GLN:OE1	2.11	0.51
1:B:43:THR:HG22	1:B:46:GLU:H	1.75	0.51
1:C:43:THR:HG22	1:C:46:GLU:H	1.76	0.51
1:G:15:GLU:OE2	1:G:91:ILE:HG23	2.11	0.51
1:H:146:TRP:HB3	1:H:147:PRO:C	2.31	0.51
1:F:115:ALA:HB3	1:F:118:LEU:HD21	1.93	0.51
1:A:122:LEU:HA	1:A:125:GLU:HB2	1.93	0.51
1:B:92:CYS:HB2	1:B:93:PRO:HD2	1.93	0.51
1:G:110:GLU:CD	1:G:110:GLU:N	2.61	0.51
1:D:134:GLN:HG2	1:D:136:TYR:HE2	1.76	0.50
1:F:19:LEU:HD22	1:F:58:ILE:HD11	1.93	0.50
1:A:63:GLN:HE22	1:A:87:GLU:H	1.58	0.50
1:F:71:TRP:CE2	1:F:122:LEU:HD13	2.47	0.50
1:D:25:ILE:HG22	1:D:26:ASN:N	2.27	0.50
1:E:28:LYS:HE2	1:E:30:LEU:HD11	1.93	0.50
1:H:31:TRP:HD1	1:H:117:ASN:ND2	2.10	0.50
1:A:65:PHE:HD2	1:A:140:MET:HE1	1.73	0.50
1:F:45:VAL:HG13	2:F:174:HOH:O	2.11	0.50
1:F:117:ASN:N	1:F:117:ASN:HD22	2.06	0.49
1:H:25:ILE:CG2	1:H:30:LEU:CD2	2.90	0.49
1:E:39:GLU:O	1:E:42:GLU:HG2	2.12	0.49
1:B:53:TRP:CZ2	1:B:97:HIS:CE1	3.00	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:81:ARG:HD2	1:H:122:LEU:HD13	1.94	0.49
1:B:63:GLN:HE22	1:B:87:GLU:HB2	1.77	0.49
1:G:38:LEU:HD13	1:G:44:LEU:HD23	1.92	0.49
1:H:22:GLU:HG3	1:H:31:TRP:CE2	2.48	0.49
1:H:71:TRP:CD1	1:H:122:LEU:HD11	2.47	0.49
1:B:36:GLY:CA	1:B:51:GLU:HG3	2.25	0.49
1:E:63:GLN:HE22	1:E:87:GLU:H	1.61	0.49
1:E:143:ASP:OD2	1:E:150:LYS:NZ	2.40	0.49
1:D:95:GLN:HG2	1:D:96:PRO:HD2	1.94	0.48
1:H:120:SER:HB2	1:H:121:PRO:HD2	1.95	0.48
1:H:31:TRP:CD1	1:H:117:ASN:ND2	2.81	0.48
1:F:22:GLU:OE2	1:F:29:ALA:HB1	2.13	0.48
1:F:66:ILE:HD11	1:F:85:ALA:HB2	1.95	0.48
1:H:20:VAL:O	1:H:105:ARG:N	2.45	0.48
1:E:31:TRP:CZ3	1:E:105:ARG:HG2	2.48	0.48
1:A:25:ILE:CG2	1:A:26:ASN:N	2.53	0.47
1:A:71:TRP:CZ3	1:A:73:ALA:HA	2.49	0.47
1:A:107:VAL:HG22	1:A:111:GLU:CB	2.44	0.47
1:F:92:CYS:HB2	1:F:93:PRO:HD2	1.96	0.47
1:C:152:VAL:HG13	1:C:153:ILE:HB	1.96	0.47
1:D:63:GLN:NE2	1:D:87:GLU:H	2.13	0.47
1:E:45:VAL:HG13	2:E:161:HOH:O	2.14	0.47
1:D:140:MET:HE3	1:F:65:PHE:CE2	2.50	0.47
1:B:136:TYR:HB3	1:B:140:MET:CE	2.43	0.47
1:A:56:THR:O	1:A:58:ILE:N	2.48	0.47
1:E:22:GLU:OE2	1:E:29:ALA:HB1	2.14	0.47
1:A:5:HIS:O	1:A:79:PHE:HA	2.14	0.47
1:B:11:VAL:HA	2:B:165:HOH:O	2.15	0.46
1:D:136:TYR:HB3	1:D:137:PRO:CD	2.45	0.46
1:D:71:TRP:CZ3	1:D:73:ALA:HA	2.51	0.46
1:E:152:VAL:O	1:F:149:THR:HG22	2.16	0.46
1:D:140:MET:HE3	1:F:65:PHE:HE2	1.81	0.46
1:A:71:TRP:CD1	1:A:122:LEU:HD11	2.51	0.46
1:F:45:VAL:HB	1:F:62:PRO:HG2	1.98	0.46
1:E:105:ARG:NH2	1:G:102:ASP:HB3	2.30	0.46
1:H:146:TRP:HB3	1:H:147:PRO:O	2.15	0.46
1:A:1:MET:HG2	1:A:3:LYS:HE2	1.97	0.46
1:E:22:GLU:HG2	1:E:102:ASP:HB2	1.98	0.46
1:G:24:THR:HA	1:G:28:LYS:O	2.15	0.46
1:A:11:VAL:HG23	1:A:83:LEU:HD11	1.97	0.46
1:D:12:VAL:HB	1:D:19:LEU:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:114:GLN:HG3	2:C:173:HOH:O	2.16	0.45
1:B:122:LEU:HA	1:B:125:GLU:HB3	1.97	0.45
1:H:71:TRP:CD1	1:H:79:PHE:HB2	2.51	0.45
1:A:72:ILE:O	1:A:72:ILE:HG13	2.15	0.45
1:D:43:THR:HG22	1:D:46:GLU:CB	2.46	0.45
1:F:134:GLN:O	1:F:134:GLN:CG	2.64	0.45
1:A:36:GLY:HA3	1:A:51:GLU:CD	2.36	0.45
1:A:140:MET:HG3	1:B:136:TYR:CE2	2.50	0.45
1:D:23:GLU:HG2	1:D:101:ILE:HD13	1.99	0.45
1:F:76:LYS:HB3	1:F:76:LYS:NZ	2.31	0.45
1:F:94:THR:HG23	1:F:96:PRO:HD3	1.98	0.45
1:E:95:GLN:HB2	1:G:93:PRO:HA	1.99	0.45
1:D:63:GLN:HE22	1:D:87:GLU:H	1.65	0.45
1:F:120:SER:O	1:F:123:VAL:HG23	2.16	0.44
1:A:56:THR:C	1:A:58:ILE:N	2.71	0.44
1:F:36:GLY:HA3	1:F:51:GLU:CG	2.47	0.44
1:G:46:GLU:OE1	1:G:150:LYS:HB2	2.17	0.44
1:F:136:TYR:HB3	1:F:137:PRO:HD2	1.99	0.44
1:A:138:LEU:O	1:A:141:ILE:HD12	2.16	0.44
1:A:152:VAL:CG1	1:A:153:ILE:N	2.81	0.44
1:E:71:TRP:CD1	1:E:79:PHE:HB2	2.52	0.44
1:C:118:LEU:HD13	1:C:123:VAL:HB	2.00	0.44
1:D:115:ALA:HB1	1:D:117:ASN:ND2	2.32	0.44
1:H:12:VAL:HB	1:H:19:LEU:HB3	2.00	0.44
1:F:36:GLY:HA3	1:F:51:GLU:CD	2.38	0.44
1:H:107:VAL:CG2	1:H:111:GLU:OE1	2.60	0.44
1:E:140:MET:HE3	1:E:140:MET:HB2	1.78	0.43
1:H:55:GLU:HG2	1:H:101:ILE:HD11	2.00	0.43
1:C:66:ILE:O	1:C:67:ARG:HB2	2.18	0.43
1:H:58:ILE:CD1	1:H:88:LEU:HD11	2.46	0.43
1:B:105:ARG:HE	1:B:105:ARG:HB3	1.65	0.43
1:C:132:SER:OG	1:C:134:GLN:HG3	2.18	0.43
1:G:134:GLN:C	1:G:134:GLN:OE1	2.56	0.43
1:A:108:SER:OG	1:A:110:GLU:HG2	2.18	0.43
1:B:91:ILE:HD12	1:B:106:TRP:CG	2.53	0.43
1:H:43:THR:HB	1:H:46:GLU:OE2	2.19	0.43
1:A:24:THR:HA	1:A:28:LYS:O	2.19	0.43
1:B:77:THR:HA	1:B:78:PRO:HD3	1.74	0.43
1:E:87:GLU:OE1	1:E:130:TYR:OH	2.31	0.43
1:F:26:ASN:C	1:F:28:LYS:H	2.21	0.43
1:F:87:GLU:OE1	1:F:130:TYR:OH	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:25:ILE:HG22	1:H:30:LEU:HD22	2.00	0.43
1:B:38:LEU:HD13	1:B:44:LEU:HD23	2.01	0.43
1:E:12:VAL:HB	1:E:19:LEU:HB3	2.01	0.42
1:E:50:ARG:HH11	1:E:50:ARG:HG2	1.84	0.42
1:E:94:THR:H	1:G:94:THR:H	1.67	0.42
1:H:89:GLU:OE2	2:H:171:HOH:O	2.22	0.42
1:E:53:TRP:O	1:E:57:GLY:HA2	2.19	0.42
1:F:75:ASP:C	1:F:76:LYS:HG2	2.40	0.42
1:B:120:SER:HB2	1:B:121:PRO:HD2	2.00	0.42
1:A:61:GLN:HE21	1:A:61:GLN:HA	1.85	0.42
1:E:130:TYR:C	1:E:130:TYR:CD2	2.93	0.42
1:C:71:TRP:NE1	1:C:122:LEU:HD11	2.34	0.42
1:E:53:TRP:O	1:E:57:GLY:N	2.53	0.41
1:G:128:ARG:NH2	1:G:128:ARG:CG	2.83	0.41
1:G:20:VAL:CG1	1:G:31:TRP:HB3	2.50	0.41
1:B:96:PRO:O	1:B:97:HIS:HB2	2.19	0.41
1:E:64:HIS:HE1	1:H:95:GLN:HE22	1.68	0.41
1:A:83:LEU:HD21	1:A:126:SER:HB3	2.01	0.41
1:C:50:ARG:NH1	1:C:51:GLU:OE2	2.53	0.41
1:D:102:ASP:O	1:D:103:CYS:HB3	2.20	0.41
1:B:54:GLU:O	1:B:96:PRO:O	2.38	0.41
1:D:11:VAL:HG11	1:D:130:TYR:CD1	2.56	0.41
1:E:96:PRO:HB3	1:E:101:ILE:HB	2.03	0.41
1:B:102:ASP:O	1:B:103:CYS:HB3	2.20	0.41
1:C:130:TYR:CD1	1:C:135:ARG:NH2	2.89	0.41
1:D:110:GLU:O	1:D:114:GLN:HG2	2.20	0.41
1:G:14:ALA:HB2	1:G:88:LEU:HB2	2.01	0.41
1:G:94:THR:HG21	1:G:104:CYS:HB2	2.02	0.41
1:G:43:THR:HG23	1:G:45:VAL:H	1.86	0.41
1:A:152:VAL:HG12	1:A:153:ILE:HG23	2.02	0.41
1:C:23:GLU:HG3	1:C:100:ASP:O	2.21	0.41
1:E:43:THR:CG2	1:E:46:GLU:H	2.25	0.41
1:G:21:VAL:HB	1:G:101:ILE:HG23	2.01	0.41
1:B:108:SER:OG	1:B:111:GLU:HG3	2.21	0.41
1:B:40:ALA:O	1:B:41:ASP:C	2.58	0.40
1:E:11:VAL:HA	2:E:156:HOH:O	2.21	0.40
1:A:118:LEU:HD22	1:A:123:VAL:HG11	2.02	0.40
1:A:3:LYS:HA	1:A:4:PRO:HD3	1.96	0.40
1:E:66:ILE:HD11	1:E:85:ALA:HB2	2.02	0.40
1:A:26:ASN:N	1:A:26:ASN:ND2	2.69	0.40
1:B:56:THR:OG1	1:B:58:ILE:HG23	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:92:CYS:HB2	1:H:106:TRP:HZ2	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	151/153 (99%)	139 (92%)	8 (5%)	4 (3%)	5	13
1	B	145/153 (95%)	135 (93%)	7 (5%)	3 (2%)	7	18
1	C	151/153 (99%)	137 (91%)	13 (9%)	1 (1%)	22	46
1	D	146/153 (95%)	132 (90%)	11 (8%)	3 (2%)	7	18
1	E	150/153 (98%)	134 (89%)	14 (9%)	2 (1%)	12	30
1	F	147/153 (96%)	134 (91%)	11 (8%)	2 (1%)	11	28
1	G	150/153 (98%)	134 (89%)	13 (9%)	3 (2%)	7	19
1	H	145/153 (95%)	131 (90%)	12 (8%)	2 (1%)	11	28
All	All	1185/1224 (97%)	1076 (91%)	89 (8%)	20 (2%)	9	23

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	100	ASP
1	H	27	GLY
1	A	57	GLY
1	C	75	ASP
1	F	41	ASP
1	F	67	ARG
1	G	103	CYS
1	B	67	ARG
1	E	26	ASN

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Mol	Chain	Res	Type
1	G	96	PRO
1	H	41	ASP
1	A	15	GLU
1	A	67	ARG
1	D	67	ARG
1	D	147	PRO
1	E	75	ASP
1	B	41	ASP
1	D	103	CYS
1	A	34	PRO
1	B	25	ILE

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	132/132 (100%)	119 (90%)	13 (10%)	8	18
1	B	127/132 (96%)	108 (85%)	19 (15%)	3	7
1	C	132/132 (100%)	116 (88%)	16 (12%)	5	11
1	D	128/132 (97%)	114 (89%)	14 (11%)	6	14
1	E	131/132 (99%)	119 (91%)	12 (9%)	9	21
1	F	129/132 (98%)	117 (91%)	12 (9%)	9	21
1	G	131/132 (99%)	110 (84%)	21 (16%)	2	6
1	H	127/132 (96%)	105 (83%)	22 (17%)	2	5
All	All	1037/1056 (98%)	908 (88%)	129 (12%)	4	11

All (129) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	VAL
1	A	26	ASN
1	A	30	LEU
1	A	52	LEU

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Mol	Chain	Res	Type
1	A	59	SER
1	A	61	GLN
1	A	72	ILE
1	A	89	GLU
1	A	105	ARG
1	A	113	LEU
1	A	127	ILE
1	A	140	MET
1	A	141	ILE
1	B	1	MET
1	B	20	VAL
1	B	30	LEU
1	B	41	ASP
1	B	43	THR
1	B	58	ILE
1	B	67	ARG
1	B	76	LYS
1	B	80	LEU
1	B	89	GLU
1	B	91	ILE
1	B	98	ASP
1	B	99	SER
1	B	107	VAL
1	B	114	GLN
1	B	116	SER
1	B	117	ASN
1	B	119	ARG
1	B	140	MET
1	C	20	VAL
1	C	30	LEU
1	C	43	THR
1	C	52	LEU
1	C	58	ILE
1	C	80	LEU
1	C	90	GLN
1	C	91	ILE
1	C	105	ARG
1	C	107	VAL
1	C	110	GLU
1	C	113	LEU
1	C	114	GLN
1	C	140	MET

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Mol	Chain	Res	Type
1	C	152	VAL
1	C	153	ILE
1	D	43	THR
1	D	59	SER
1	D	67	ARG
1	D	76	LYS
1	D	80	LEU
1	D	94	THR
1	D	95	GLN
1	D	105	ARG
1	D	107	VAL
1	D	117	ASN
1	D	125	GLU
1	D	127	ILE
1	D	134	GLN
1	D	140	MET
1	E	3	LYS
1	E	20	VAL
1	E	26	ASN
1	E	30	LEU
1	E	43	THR
1	E	52	LEU
1	E	58	ILE
1	E	67	ARG
1	E	76	LYS
1	E	80	LEU
1	E	94	THR
1	E	104	CYS
1	F	1	MET
1	F	26	ASN
1	F	30	LEU
1	F	43	THR
1	F	67	ARG
1	F	76	LYS
1	F	89	GLU
1	F	94	THR
1	F	117	ASN
1	F	122	LEU
1	F	132	SER
1	F	149	THR
1	G	25	ILE
1	G	42	GLU

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Mol	Chain	Res	Type
1	G	43	THR
1	G	59	SER
1	G	67	ARG
1	G	76	LYS
1	G	77	THR
1	G	80	LEU
1	G	89	GLU
1	G	90	GLN
1	G	91	ILE
1	G	94	THR
1	G	95	GLN
1	G	98	ASP
1	G	107	VAL
1	G	110	GLU
1	G	116	SER
1	G	134	GLN
1	G	135	ARG
1	G	149	THR
1	G	150	LYS
1	H	1	MET
1	H	17	LYS
1	H	26	ASN
1	H	30	LEU
1	H	41	ASP
1	H	42	GLU
1	H	43	THR
1	H	52	LEU
1	H	58	ILE
1	H	76	LYS
1	H	77	THR
1	H	80	LEU
1	H	89	GLU
1	H	92	CYS
1	H	94	THR
1	H	100	ASP
1	H	102	ASP
1	H	105	ARG
1	H	114	GLN
1	H	119	ARG
1	H	128	ARG
1	H	146	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	HIS
1	A	26	ASN
1	A	61	GLN
1	A	63	GLN
1	B	63	GLN
1	B	117	ASN
1	B	145	ASN
1	C	63	GLN
1	C	114	GLN
1	D	26	ASN
1	D	63	GLN
1	D	90	GLN
1	D	95	GLN
1	D	117	ASN
1	D	134	GLN
1	D	145	ASN
1	E	26	ASN
1	E	61	GLN
1	E	63	GLN
1	E	95	GLN
1	F	5	HIS
1	F	37	HIS
1	F	61	GLN
1	F	63	GLN
1	F	117	ASN
1	G	61	GLN
1	G	63	GLN
1	G	95	GLN
1	G	114	GLN
1	H	63	GLN
1	H	95	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 [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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	153/153 (100%)	-0.41	1 (0%) 87 89	15, 29, 44, 51	2 (1%)
1	B	147/153 (96%)	-0.43	0 100 100	15, 26, 39, 56	0
1	C	153/153 (100%)	-0.31	2 (1%) 77 78	13, 27, 39, 44	0
1	D	148/153 (96%)	-0.68	0 100 100	13, 24, 35, 44	2 (1%)
1	E	152/153 (99%)	-0.44	0 100 100	15, 26, 40, 53	0
1	F	149/153 (97%)	-0.49	0 100 100	11, 27, 42, 53	1 (0%)
1	G	152/153 (99%)	-0.45	0 100 100	15, 29, 45, 55	0
1	H	147/153 (96%)	-0.49	1 (0%) 87 89	15, 27, 45, 62	0
All	All	1201/1224 (98%)	-0.46	4 (0%) 94 95	11, 27, 43, 62	5 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1	MET	3.2
1	H	146	TRP	3.1
1	C	115	ALA	2.2
1	A	139	GLU	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 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.