



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 02:27 am GMT

PDB ID : 1VDM  
Title : Crystal structure of purine phosphoribosyltransferase from *Pyrococcus horikoshii* Ot3  
Authors : Sugahara, M.; Kunishima, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 2004-03-23  
Resolution : 2.50 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

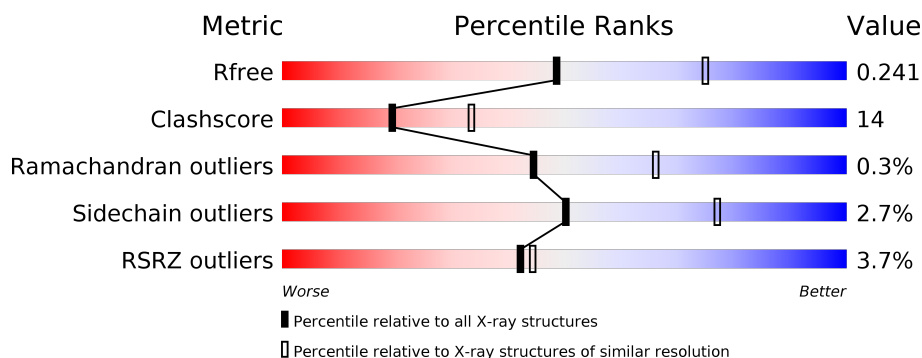
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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	153	<div> <div>7%</div> <div> <div></div> <div>62%</div> <div>34%</div> <div>...</div> </div> </div>
1	B	153	<div> <div>3%</div> <div> <div></div> <div>64%</div> <div>33%</div> <div>..</div> </div> </div>
1	C	153	<div> <div>5%</div> <div> <div></div> <div>69%</div> <div>24%</div> <div>7%</div> </div> </div>
1	D	153	<div> <div>6%</div> <div> <div></div> <div>68%</div> <div>27%</div> <div>..</div> </div> </div>
1	E	153	<div> <div>5%</div> <div> <div></div> <div>66%</div> <div>30%</div> <div>..</div> </div> </div>
1	F	153	<div> <div></div> <div> <div></div> <div>82%</div> <div>15%</div> <div>..</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	153	<div><div><div></div><div></div><div></div></div><div>2%64%33%<div><div></div><div></div></div></div></div>
1	H	153	<div><div><div></div><div></div><div></div></div><div>4%64%31%<div><div></div><div></div></div></div></div>
1	I	153	<div><div><div></div><div></div><div></div></div><div>%71%24%<div><div></div><div></div></div></div></div>
1	J	153	<div><div><div></div><div></div><div></div></div><div>%75%20%<div><div></div><div></div></div></div></div>
1	K	153	<div><div><div></div><div></div><div></div></div><div>5%61%33%<div><div></div><div></div></div></div></div>
1	L	153	<div><div><div></div><div></div><div></div></div><div>5%67%30%<div><div></div><div></div></div></div></div>

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 15181 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 purine phosphoribosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	151	Total	C	N	O	S	0	0	0
			1227	810	199	215	3			
1	B	149	Total	C	N	O	S	0	0	0
			1207	799	194	211	3			
1	C	143	Total	C	N	O	S	0	0	0
			1159	770	187	199	3			
1	D	147	Total	C	N	O	S	0	0	0
			1193	792	192	206	3			
1	E	149	Total	C	N	O	S	0	0	0
			1211	803	194	211	3			
1	F	150	Total	C	N	O	S	0	0	0
			1213	804	195	211	3			
1	G	152	Total	C	N	O	S	0	0	0
			1233	815	200	215	3			
1	H	149	Total	C	N	O	S	0	0	0
			1214	803	197	211	3			
1	I	146	Total	C	N	O	S	0	0	0
			1184	786	190	205	3			
1	J	147	Total	C	N	O	S	0	0	0
			1195	792	194	206	3			
1	K	146	Total	C	N	O	S	0	0	0
			1189	790	191	205	3			
1	L	149	Total	C	N	O	S	0	0	0
			1207	799	194	211	3			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	46	Total	O	0	0
			46	46		
2	B	69	Total	O	0	0
			69	69		

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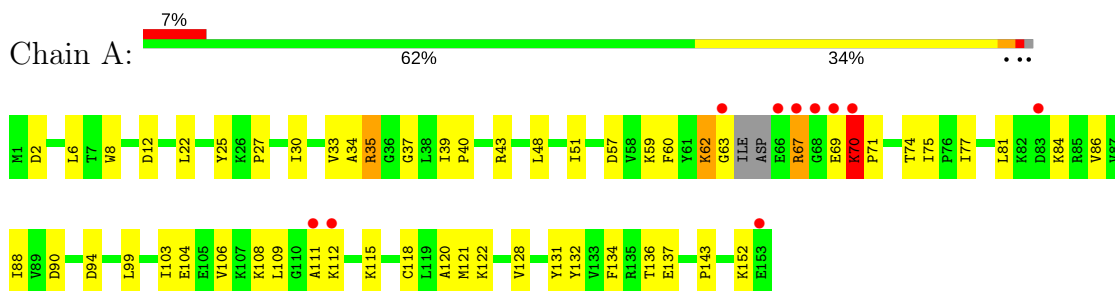
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	52	Total 52	O 52	0	0
2	D	57	Total 57	O 57	0	0
2	E	92	Total 92	O 92	0	0
2	F	98	Total 98	O 98	0	0
2	G	52	Total 52	O 52	0	0
2	H	51	Total 51	O 51	0	0
2	I	52	Total 52	O 52	0	0
2	J	59	Total 59	O 59	0	0
2	K	52	Total 52	O 52	0	0
2	L	69	Total 69	O 69	0	0

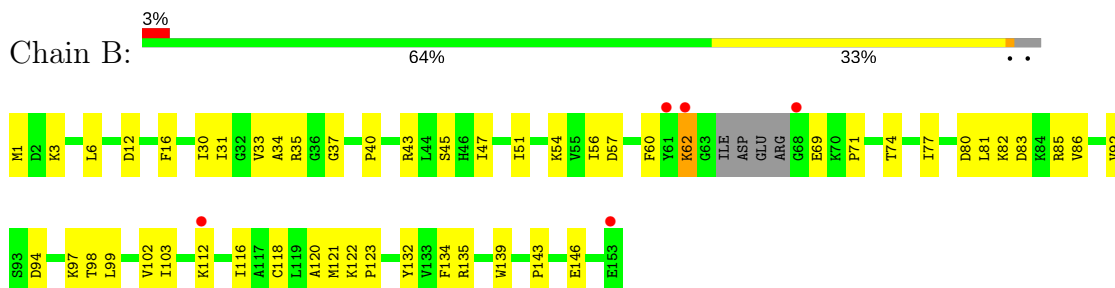
### 3 Residue-property plots [i](#)

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.

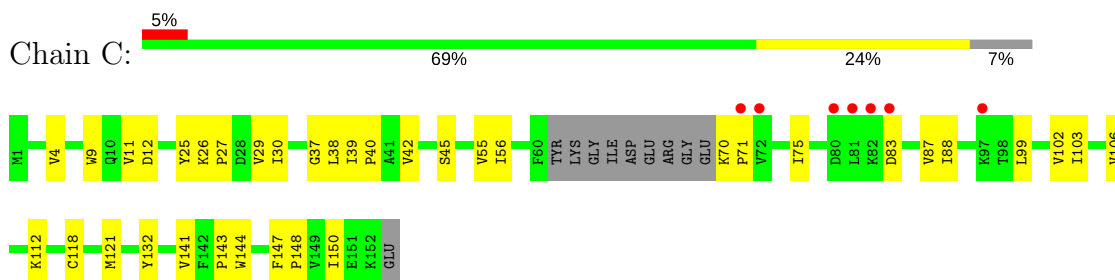
- Molecule 1: purine phosphoribosyltransferase



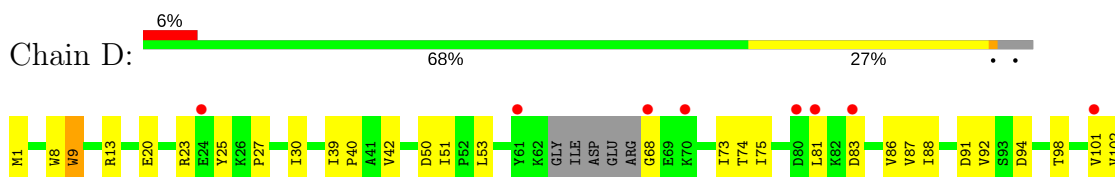
- Molecule 1: purine phosphoribosyltransferase

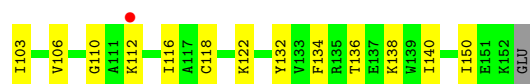


- Molecule 1: purine phosphoribosyltransferase

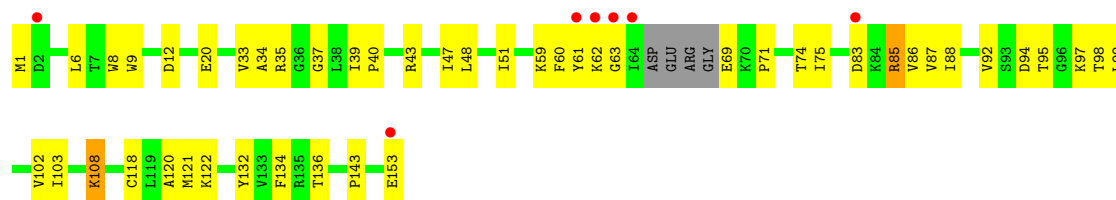


- Molecule 1: purine phosphoribosyltransferase

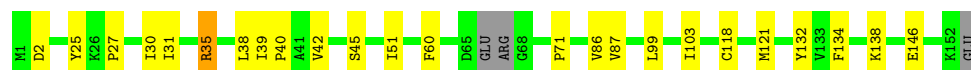
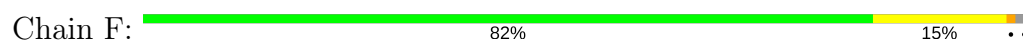




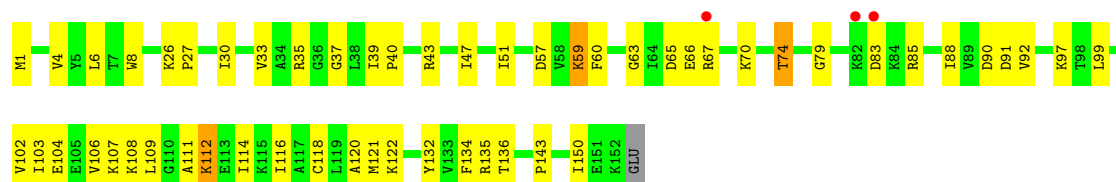
- Molecule 1: purine phosphoribosyltransferase



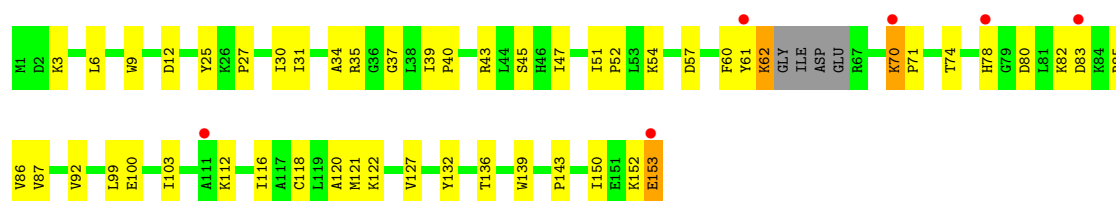
- Molecule 1: purine phosphoribosyltransferase



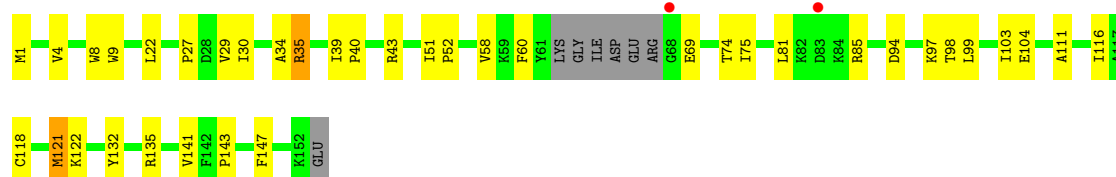
- Molecule 1: purine phosphoribosyltransferase



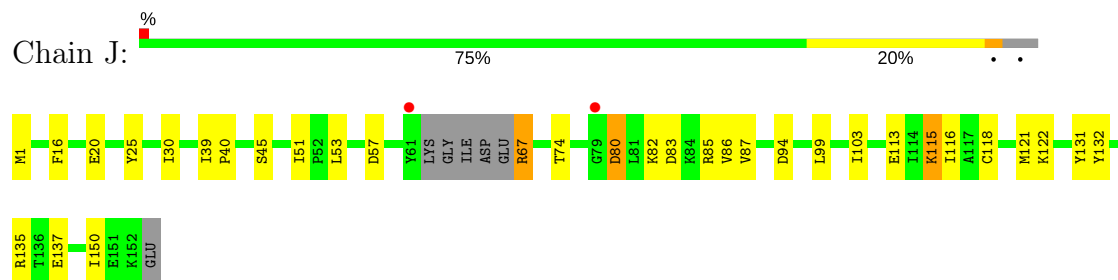
- Molecule 1: purine phosphoribosyltransferase



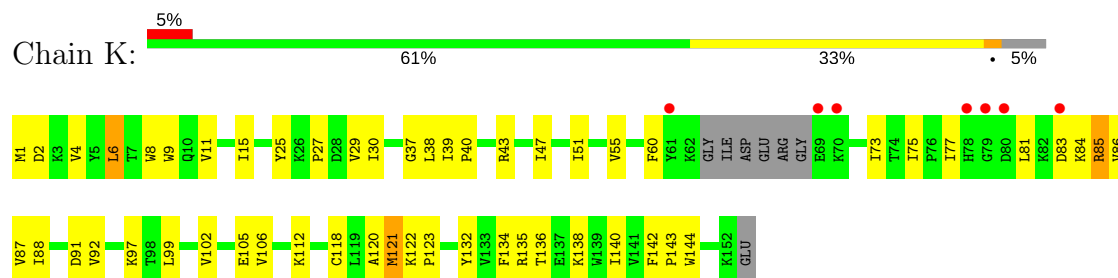
- Molecule 1: purine phosphoribosyltransferase



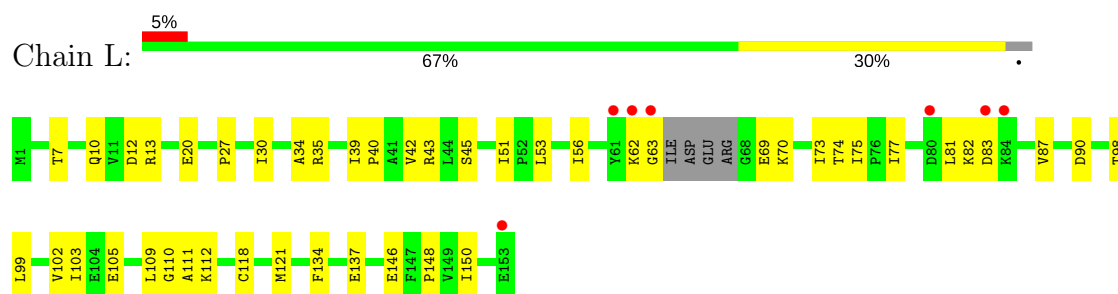
## • Molecule 1: purine phosphoribosyltransferase



## • Molecule 1: purine phosphoribosyltransferase



## • Molecule 1: purine phosphoribosyltransferase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	105.60Å 157.35Å 170.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.55 – 2.50 39.55 – 2.49	Depositor EDS
% Data completeness (in resolution range)	100.0 (39.55-2.50) 99.7 (39.55-2.49)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.50 (at 2.48Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.209 , 0.241 0.209 , 0.241	Depositor DCC
$R_{free}$ test set	4938 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	42.0	Xtriage
Anisotropy	0.454	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 57.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	15181	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.67% 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

### 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	0.44	0/1256	0.68	0/1701
1	B	0.44	0/1236	0.63	0/1675
1	C	0.39	0/1187	0.64	0/1612
1	D	0.54	2/1222 (0.2%)	0.63	0/1658
1	E	0.43	0/1240	0.65	0/1681
1	F	0.41	0/1242	0.64	0/1685
1	G	0.43	0/1263	0.67	0/1714
1	H	0.42	0/1243	0.64	0/1684
1	I	0.40	0/1213	0.63	0/1647
1	J	0.42	0/1224	0.62	0/1661
1	K	0.46	0/1218	0.67	0/1653
1	L	0.40	0/1236	0.59	0/1675
All	All	0.43	2/14780 (0.0%)	0.64	0/20046

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	9	TRP	NE1-CE2	8.76	1.49	1.37
1	D	8	TRP	NE1-CE2	8.73	1.49	1.37

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	1284	47	0
1	B	1207	0	1265	36	0
1	C	1159	0	1225	36	0
1	D	1193	0	1256	38	0
1	E	1211	0	1273	40	0
1	F	1213	0	1274	19	0
1	G	1233	0	1294	51	0
1	H	1214	0	1275	44	0
1	I	1184	0	1243	33	0
1	J	1195	0	1256	25	0
1	K	1189	0	1253	54	0
1	L	1207	0	1265	39	0
2	A	46	0	0	0	0
2	B	69	0	0	1	0
2	C	52	0	0	0	0
2	D	57	0	0	1	0
2	E	92	0	0	2	0
2	F	98	0	0	0	0
2	G	52	0	0	1	0
2	H	51	0	0	0	0
2	I	52	0	0	0	0
2	J	59	0	0	1	0
2	K	52	0	0	1	0
2	L	69	0	0	0	0
All	All	15181	0	15163	425	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 14.

The worst 5 of 425 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:70:LYS:HB3	1:A:71:PRO:HD2	1.28	1.07
1:C:70:LYS:HB2	1:C:71:PRO:HD3	1.43	1.00
1:J:30:ILE:HG13	1:J:51:ILE:HD11	1.45	0.97
1:G:66:GLU:HG2	1:G:67:ARG:N	1.82	0.94
1:C:39:ILE:HB	1:C:40:PRO:HD3	1.49	0.92

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	147/153 (96%)	135 (92%)	10 (7%)	2 (1%)	13	23
1	B	145/153 (95%)	139 (96%)	6 (4%)	0	100	100
1	C	139/153 (91%)	129 (93%)	10 (7%)	0	100	100
1	D	143/153 (94%)	133 (93%)	8 (6%)	2 (1%)	13	23
1	E	145/153 (95%)	138 (95%)	7 (5%)	0	100	100
1	F	146/153 (95%)	143 (98%)	3 (2%)	0	100	100
1	G	150/153 (98%)	144 (96%)	5 (3%)	1 (1%)	25	43
1	H	145/153 (95%)	139 (96%)	6 (4%)	0	100	100
1	I	142/153 (93%)	135 (95%)	7 (5%)	0	100	100
1	J	143/153 (94%)	138 (96%)	5 (4%)	0	100	100
1	K	142/153 (93%)	135 (95%)	7 (5%)	0	100	100
1	L	145/153 (95%)	140 (97%)	5 (3%)	0	100	100
All	All	1732/1836 (94%)	1648 (95%)	79 (5%)	5 (0%)	44	66

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	152	LYS
1	G	112	LYS
1	A	70	LYS
1	D	110	GLY
1	D	73	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	134/136 (98%)	128 (96%)	6 (4%)	32	56
1	B	132/136 (97%)	130 (98%)	2 (2%)	70	89
1	C	128/136 (94%)	126 (98%)	2 (2%)	68	88
1	D	131/136 (96%)	129 (98%)	2 (2%)	70	89
1	E	133/136 (98%)	129 (97%)	4 (3%)	46	74
1	F	133/136 (98%)	131 (98%)	2 (2%)	70	89
1	G	135/136 (99%)	130 (96%)	5 (4%)	39	66
1	H	133/136 (98%)	128 (96%)	5 (4%)	38	64
1	I	130/136 (96%)	127 (98%)	3 (2%)	56	81
1	J	131/136 (96%)	125 (95%)	6 (5%)	31	55
1	K	131/136 (96%)	128 (98%)	3 (2%)	56	81
1	L	132/136 (97%)	130 (98%)	2 (2%)	70	89
All	All	1583/1632 (97%)	1541 (97%)	42 (3%)	50	77

5 of 42 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	35	ARG
1	H	62	LYS
1	K	85	ARG
1	G	59	LYS
1	G	74	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	HIS

### 5.3.3 RNA ⓘ

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 [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	151/153 (98%)	0.12	10 (6%)	19	19	24, 46, 75, 104	0
1	B	149/153 (97%)	0.08	5 (3%)	46	48	23, 41, 69, 91	0
1	C	143/153 (93%)	0.25	7 (4%)	30	32	24, 51, 81, 85	0
1	D	147/153 (96%)	0.26	9 (6%)	22	22	25, 48, 87, 97	0
1	E	149/153 (97%)	0.05	7 (4%)	32	34	23, 40, 68, 88	0
1	F	150/153 (98%)	-0.13	0	100	100	22, 38, 59, 82	0
1	G	152/153 (99%)	0.13	3 (1%)	65	67	23, 48, 78, 89	0
1	H	149/153 (97%)	0.05	6 (4%)	39	41	23, 44, 80, 86	0
1	I	146/153 (95%)	-0.05	2 (1%)	75	76	20, 42, 64, 79	0
1	J	147/153 (96%)	-0.19	2 (1%)	75	76	22, 37, 62, 75	0
1	K	146/153 (95%)	0.08	7 (4%)	31	32	23, 40, 73, 84	0
1	L	149/153 (97%)	0.10	7 (4%)	32	34	23, 46, 79, 89	0
All	All	1778/1836 (96%)	0.06	65 (3%)	42	44	20, 44, 78, 104	0

The worst 5 of 65 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	78	HIS	4.5
1	I	68	GLY	4.2
1	E	153	GLU	4.2
1	A	66	GLU	4.1
1	B	153	GLU	4.0

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

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

### 6.3 Carbohydrates

There are no carbohydrates in this entry.

### 6.4 Ligands

There are no ligands in this entry.

### 6.5 Other polymers

There are no such residues in this entry.