



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:23 AM GMT

PDB ID : 2GVL  
Title : Crystal Structure of Murine NMPRTase  
Authors : Khan, J.A.; Tao, X.; Tong, L.  
Deposited on : 2006-05-02  
Resolution : 2.10 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

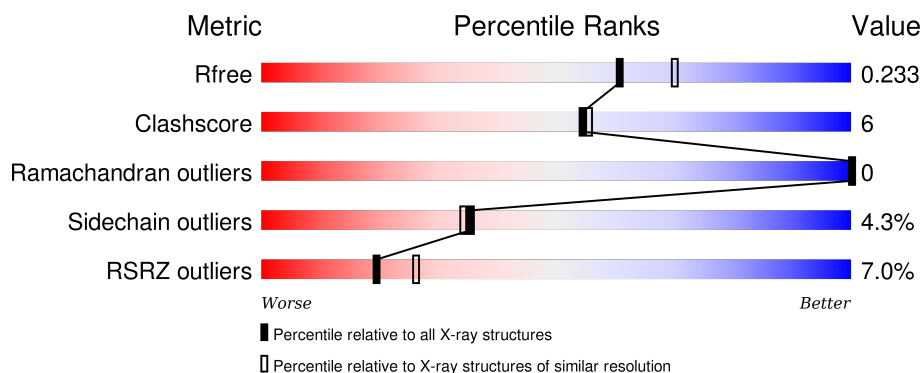
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


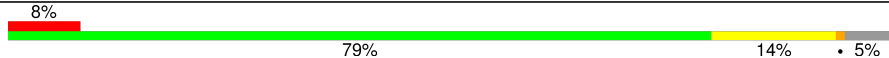
The reported resolution of this entry is 2.10 Å.

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))
$R_{free}$	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	491	
1	B	491	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8133 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 Nicotinamide phosphoribosyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	Se	0	0	0
			3710	2381	618	705	5	1			
1	B	464	Total	C	N	O	S	Se	0	0	0
			3710	2381	618	705	5	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	368	MSE	MET	MODIFIED RESIDUE	UNP Q99KQ4
B	368	MSE	MET	MODIFIED RESIDUE	UNP Q99KQ4

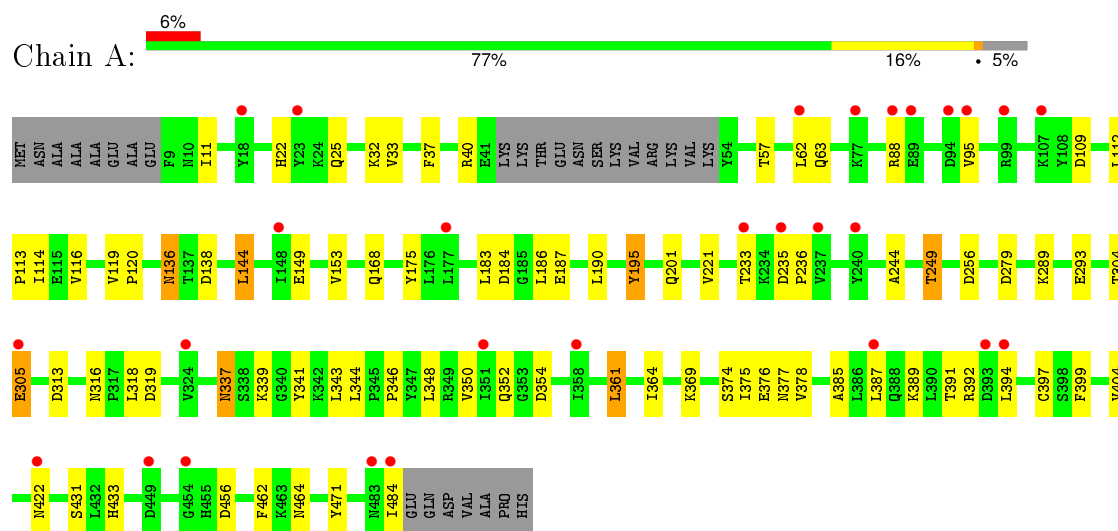
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	348	Total	O	0	0
			348	348		
2	B	365	Total	O	0	0
			365	365		

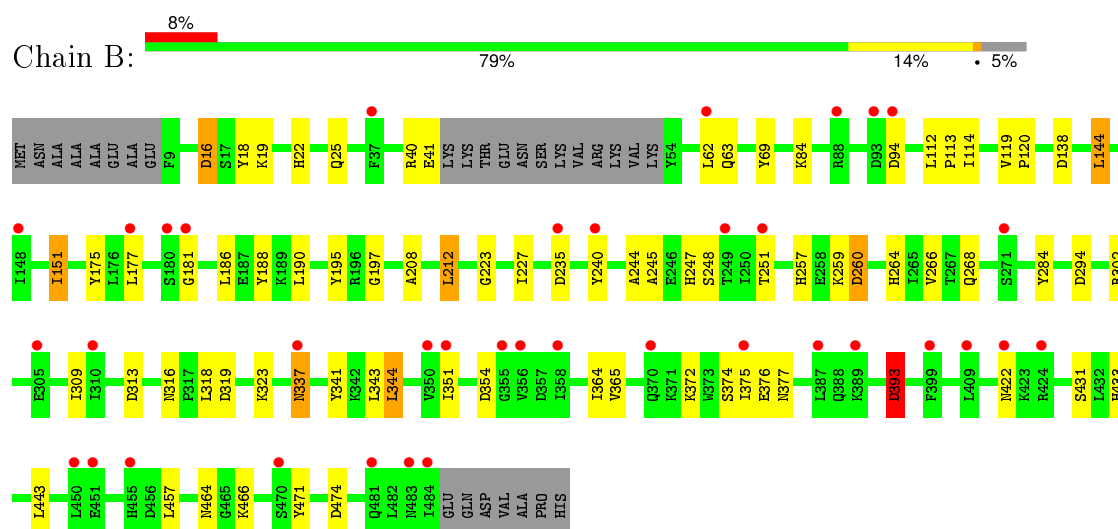
### 3 Residue-property plots

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: Nicotinamide phosphoribosyltransferase



- Molecule 1: Nicotinamide phosphoribosyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	60.26 Å 107.74 Å 83.29 Å 90.00° 96.57° 90.00°	Depositor
Resolution (Å)	82.76 – 2.10 29.94 – 2.10	Depositor EDS
% Data completeness (in resolution range)	94.8 (82.76-2.10) 90.6 (29.94-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.86 (at 2.10 Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, $R_{free}$	0.225 , 0.270 0.227 , 0.233	Depositor DCC
$R_{free}$ test set	4491 reflections (8.69%)	DCC
Wilson B-factor (Å <sup>2</sup> )	18.9	Xtriage
Anisotropy	0.510	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 49.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 58582 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	8133	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.98% 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

### 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	1/3797 (0.0%)	0.70	7/5147 (0.1%)
1	B	0.43	0/3797	0.71	8/5147 (0.2%)
All	All	0.43	1/7594 (0.0%)	0.70	15/10294 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	305	GLU	CD-OE2	7.13	1.33	1.25

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	456	ASP	CB-CG-OD2	7.66	125.19	118.30
1	B	138	ASP	CB-CG-OD2	6.31	123.98	118.30
1	A	138	ASP	CB-CG-OD2	6.10	123.79	118.30
1	B	313	ASP	CB-CG-OD2	5.83	123.55	118.30
1	B	16	ASP	CB-CG-OD2	5.82	123.54	118.30
1	B	474	ASP	CB-CG-OD2	5.60	123.34	118.30
1	A	313	ASP	CB-CG-OD2	5.26	123.03	118.30
1	A	256	ASP	CB-CG-OD2	5.24	123.02	118.30
1	B	393	ASP	CB-CG-OD2	5.19	122.97	118.30
1	B	260	ASP	CB-CG-OD2	5.17	122.95	118.30
1	B	354	ASP	CB-CG-OD2	5.17	122.95	118.30
1	B	235	ASP	CB-CG-OD2	5.13	122.91	118.30
1	A	109	ASP	CB-CG-OD2	5.12	122.91	118.30
1	A	184	ASP	CB-CG-OD2	5.11	122.90	118.30
1	A	279	ASP	CB-CG-OD2	5.09	122.88	118.30

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	3710	0	3677	49	0
1	B	3710	0	3677	51	0
2	A	348	0	0	5	0
2	B	365	0	0	3	0
All	All	8133	0	7354	94	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 6.

All (94) 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:348:LEU:HB3	2:A:571:HOH:O	1.72	0.88
1:A:344:LEU:H	1:A:377:ASN:HD21	1.25	0.84
1:B:40:ARG:HD2	1:B:422:ASN:O	1.80	0.81
1:B:344:LEU:H	1:B:377:ASN:HD21	1.30	0.79
1:B:63:GLN:HE22	1:B:471:TYR:H	1.32	0.76
1:A:25:GLN:HE22	1:B:245:ALA:HA	1.53	0.74
1:A:195:TYR:OH	1:A:201:GLN:NE2	2.22	0.72
2:A:498:HOH:O	1:B:251:THR:HG21	1.90	0.71
1:A:352:GLN:HE21	1:A:354:ASP:H	1.39	0.71
1:A:344:LEU:H	1:A:377:ASN:ND2	1.90	0.70
1:B:316:ASN:HD22	1:B:319:ASP:H	1.40	0.69
1:B:365:VAL:HG13	1:B:375:ILE:HD12	1.73	0.69
1:A:249:THR:HG21	1:B:25:GLN:OE1	1.94	0.68
1:B:365:VAL:HG13	1:B:375:ILE:CD1	2.22	0.68
1:B:264:HIS:O	1:B:268:GLN:HG2	1.96	0.66
1:A:316:ASN:HD22	1:A:319:ASP:H	1.43	0.66
1:A:337:ASN:C	1:A:337:ASN:HD22	2.00	0.64
1:A:343:LEU:HD13	1:A:376:GLU:HG3	1.77	0.64
1:B:343:LEU:CD1	1:B:376:GLU:HG3	2.29	0.62
1:A:431:SER:OG	1:A:433:HIS:HE1	1.82	0.62
1:B:337:ASN:N	1:B:337:ASN:HD22	1.98	0.61
1:B:343:LEU:HD13	1:B:376:GLU:HG3	1.81	0.61
1:B:114:ILE:HG23	1:B:144:LEU:HD13	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ILE:HG12	1:A:144:LEU:HD13	1.83	0.59
1:B:344:LEU:H	1:B:377:ASN:ND2	2.00	0.58
1:B:393:ASP:N	1:B:393:ASP:OD1	2.38	0.57
1:A:119:VAL:HG22	1:A:120:PRO:HD2	1.86	0.57
1:A:385:ALA:HA	1:A:389:LYS:HB3	1.86	0.57
1:A:343:LEU:CD1	1:A:376:GLU:HG3	2.35	0.57
1:A:374:SER:OG	1:A:376:GLU:HG2	2.05	0.57
1:A:32:LYS:HA	1:A:136:ASN:HD21	1.71	0.56
1:B:208:ALA:O	1:B:212:LEU:HD22	2.05	0.56
1:A:63:GLN:HE22	1:A:471:TYR:H	1.53	0.56
1:B:181:GLY:HA3	2:B:528:HOH:O	2.06	0.55
1:B:223:GLY:O	1:B:227:ILE:HG12	2.07	0.54
1:A:37:PHE:CZ	1:A:397:CYS:HB3	2.43	0.54
1:B:337:ASN:ND2	1:B:341:TYR:O	2.41	0.54
1:A:175:TYR:HB3	1:A:375:ILE:HG13	1.91	0.52
1:A:187:GLU:HB2	2:A:768:HOH:O	2.08	0.52
1:B:113:PRO:HD2	1:B:144:LEU:CD1	2.40	0.52
1:B:431:SER:OG	1:B:433:HIS:HE1	1.93	0.52
1:A:348:LEU:CB	2:A:571:HOH:O	2.41	0.52
1:B:175:TYR:HB3	1:B:375:ILE:HG13	1.92	0.51
1:B:188:TYR:HD2	1:B:240:TYR:CE1	2.29	0.50
1:A:350:VAL:HB	1:A:378:VAL:HG22	1.93	0.50
1:A:136:ASN:HD22	1:A:136:ASN:H	1.59	0.50
1:B:318:LEU:HD13	1:B:364:ILE:HA	1.93	0.50
1:A:40:ARG:HD2	1:A:422:ASN:O	2.12	0.50
1:A:391:THR:H	1:A:394:LEU:HD13	1.76	0.50
1:A:33:VAL:H	1:A:136:ASN:ND2	2.10	0.49
1:B:247:HIS:O	1:B:251:THR:HG22	2.11	0.49
1:B:248:SER:HA	1:B:251:THR:HG22	1.94	0.49
1:A:40:ARG:NH1	1:A:422:ASN:O	2.46	0.48
1:A:113:PRO:HA	1:A:464:ASN:HD22	1.79	0.47
1:A:389:LYS:HG3	1:A:389:LYS:O	2.14	0.47
1:A:221:VAL:HG22	2:B:505:HOH:O	2.14	0.47
1:B:284:TYR:CE2	1:B:323:LYS:HD2	2.50	0.46
1:B:19:LYS:HA	1:B:22:HIS:CD2	2.51	0.46
1:A:392:ARG:HD2	1:B:197:GLY:HA2	1.98	0.46
1:B:16:ASP:HB3	1:B:18:TYR:CE2	2.51	0.46
1:B:343:LEU:HD13	1:B:376:GLU:CG	2.47	0.45
1:A:32:LYS:O	1:A:404:VAL:HG23	2.16	0.45
1:B:284:TYR:CD2	1:B:323:LYS:HD2	2.50	0.45
1:B:337:ASN:HD21	1:B:343:LEU:HB2	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:VAL:H	1:A:136:ASN:HD21	1.64	0.45
1:B:113:PRO:HD2	1:B:144:LEU:HD12	1.97	0.45
1:A:289:LYS:O	1:A:293:GLU:HB2	2.17	0.45
1:A:116:VAL:HB	1:A:462:PHE:HB3	1.99	0.45
1:B:175:TYR:CD2	1:B:365:VAL:HG12	2.52	0.43
1:B:119:VAL:HG22	1:B:120:PRO:HD2	2.00	0.43
1:A:316:ASN:ND2	1:A:319:ASP:H	2.11	0.43
1:B:374:SER:OG	1:B:376:GLU:HG2	2.18	0.43
1:A:392:ARG:CD	1:B:197:GLY:HA2	2.49	0.43
1:A:318:LEU:HD13	1:A:364:ILE:HA	2.00	0.43
1:B:266:VAL:O	1:B:302:ARG:NH2	2.52	0.42
1:A:304:THR:HG23	1:A:346:PRO:HB2	2.02	0.42
1:B:257:HIS:HD2	1:B:260:ASP:OD2	2.03	0.42
1:A:387:LEU:HD11	2:A:517:HOH:O	2.20	0.42
1:B:309:ILE:HG22	1:B:351:ILE:HG22	2.01	0.42
1:A:22:HIS:HE1	1:B:244:ALA:O	2.03	0.41
1:A:337:ASN:ND2	1:A:341:TYR:H	2.19	0.41
1:B:259:LYS:HD2	1:B:294:ASP:HB3	2.02	0.41
1:B:337:ASN:N	1:B:337:ASN:ND2	2.66	0.40
1:A:168:GLN:HG2	1:A:361:LEU:HD12	2.03	0.40
1:B:433:HIS:CE1	1:B:443:LEU:HD12	2.56	0.40
1:A:244:ALA:O	1:B:22:HIS:HE1	2.05	0.40
1:B:69:TYR:CD1	1:B:151:ILE:HD11	2.57	0.40
1:B:175:TYR:CG	1:B:365:VAL:HG12	2.56	0.40
1:A:149:GLU:HG3	1:A:399:PHE:CD2	2.57	0.40
1:A:112:LEU:HA	1:A:113:PRO:HD3	1.81	0.40
1:A:235:ASP:HA	1:A:236:PRO:HD3	1.87	0.40
1:B:84:LYS:NZ	2:B:843:HOH:O	2.53	0.40
1:A:153:VAL:CG2	1:A:399:PHE:HB2	2.51	0.40
1:B:112:LEU:O	1:B:464:ASN:HA	2.22	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	460/491 (94%)	447 (97%)	13 (3%)	0	100	100
1	B	460/491 (94%)	442 (96%)	18 (4%)	0	100	100
All	All	920/982 (94%)	889 (97%)	31 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	409/430 (95%)	390 (95%)	19 (5%)	33	31
1	B	409/430 (95%)	393 (96%)	16 (4%)	39	39
All	All	818/860 (95%)	783 (96%)	35 (4%)	35	34

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ILE
1	A	57	THR
1	A	62	LEU
1	A	88	ARG
1	A	95	VAL
1	A	136	ASN
1	A	144	LEU
1	A	183	LEU
1	A	186	LEU
1	A	190	LEU
1	A	195	TYR
1	A	233	THR
1	A	249	THR
1	A	305	GLU
1	A	337	ASN

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Mol	Chain	Res	Type
1	A	339	LYS
1	A	361	LEU
1	A	369	LYS
1	A	484	ILE
1	B	41	GLU
1	B	62	LEU
1	B	94	ASP
1	B	144	LEU
1	B	151	ILE
1	B	177	LEU
1	B	186	LEU
1	B	190	LEU
1	B	195	TYR
1	B	212	LEU
1	B	337	ASN
1	B	344	LEU
1	B	372	LYS
1	B	393	ASP
1	B	457	LEU
1	B	466	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	HIS
1	A	25	GLN
1	A	63	GLN
1	A	102	ASN
1	A	129	ASN
1	A	136	ASN
1	A	201	GLN
1	A	257	HIS
1	A	316	ASN
1	A	337	ASN
1	A	352	GLN
1	A	362	GLN
1	A	377	ASN
1	A	396	ASN
1	A	433	HIS
1	A	459	HIS
1	A	464	ASN
1	A	479	ASN

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Mol	Chain	Res	Type
1	B	22	HIS
1	B	63	GLN
1	B	92	GLN
1	B	97	ASN
1	B	146	ASN
1	B	201	GLN
1	B	264	HIS
1	B	316	ASN
1	B	337	ASN
1	B	377	ASN
1	B	396	ASN
1	B	412	ASN
1	B	433	HIS
1	B	459	HIS
1	B	479	ASN
1	B	481	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 ⓘ

### 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	463/491 (94%)	0.69	28 (6%) 25 33	10, 19, 30, 35	0
1	B	463/491 (94%)	0.74	37 (7%) 15 21	12, 20, 30, 37	0
All	All	926/982 (94%)	0.72	65 (7%) 19 26	10, 20, 30, 37	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	484	ILE	6.1
1	A	88	ARG	5.0
1	B	235	ASP	4.6
1	B	88	ARG	4.2
1	A	95	VAL	3.9
1	A	237	VAL	3.7
1	B	484	ILE	3.7
1	B	451	GLU	3.5
1	A	454	GLY	3.5
1	B	180	SER	3.3
1	A	18	TYR	3.2
1	A	235	ASP	3.2
1	B	148	ILE	3.2
1	B	271	SER	3.1
1	A	393	ASP	3.0
1	B	251	THR	3.0
1	B	483	ASN	3.0
1	B	351	ILE	2.9
1	A	351	ILE	2.8
1	B	350	VAL	2.8
1	A	305	GLU	2.8
1	B	455	HIS	2.8
1	B	370	GLN	2.7
1	B	181	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	394	LEU	2.7
1	A	89	GLU	2.7
1	B	240	TYR	2.7
1	A	148	ILE	2.6
1	A	387	LEU	2.5
1	A	94	ASP	2.5
1	A	77	LYS	2.5
1	A	358	ILE	2.5
1	B	337	ASN	2.5
1	B	409	LEU	2.5
1	B	450	LEU	2.5
1	B	310	ILE	2.5
1	A	483	ASN	2.4
1	A	62	LEU	2.4
1	A	99	ARG	2.4
1	B	424	ARG	2.4
1	B	481	GLN	2.4
1	B	389	LYS	2.4
1	B	355	GLY	2.4
1	B	358	ILE	2.4
1	B	94	ASP	2.3
1	B	387	LEU	2.3
1	A	177	LEU	2.3
1	B	177	LEU	2.3
1	A	324	VAL	2.3
1	B	422	ASN	2.3
1	A	449	ASP	2.2
1	B	62	LEU	2.2
1	B	305	GLU	2.2
1	B	399	PHE	2.2
1	B	93	ASP	2.2
1	A	422	ASN	2.2
1	A	233	THR	2.1
1	A	23	TYR	2.1
1	B	470	SER	2.1
1	A	240	TYR	2.1
1	B	356	VAL	2.1
1	B	249	THR	2.0
1	A	107	LYS	2.0
1	B	375	ILE	2.0
1	B	37	PHE	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.