



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:03 PM GMT

PDB ID : 1WD7  
Title : Crystal Structure of Uroporphyrinogen III Synthase from an Extremely Thermophilic Bacterium *Thermus thermophilus* HB8 (Wild type, Native, Form-2 crystal)  
Authors : Mizohata, E.; Matsuura, T.; Sakai, H.; Murayama, K.; Terada, T.; Shirouzu, M.; Kuramitsu, S.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 2004-05-12  
Resolution : 1.60 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

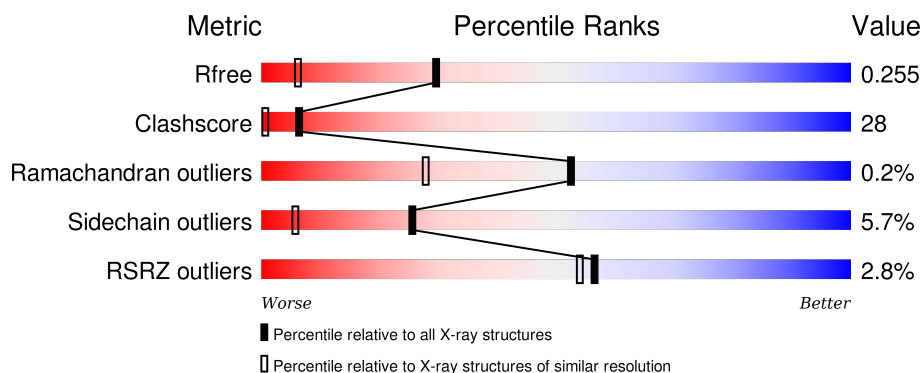
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

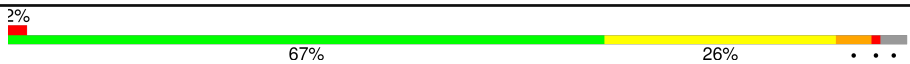
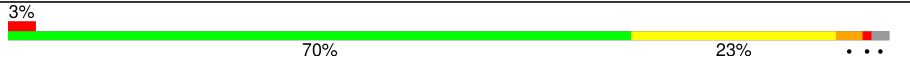
The reported resolution of this entry is 1.60 Å.

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	2475 (1.60-1.60)
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)
RSRZ outliers	91569	2479 (1.60-1.60)

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	261	
1	B	261	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4357 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 Uroporphyrinogen III Synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	254	Total	C	N	O	S	0	0	0
			1943	1252	348	342	1			
1	B	255	Total	C	N	O	S	0	0	0
			1951	1256	349	345	1			

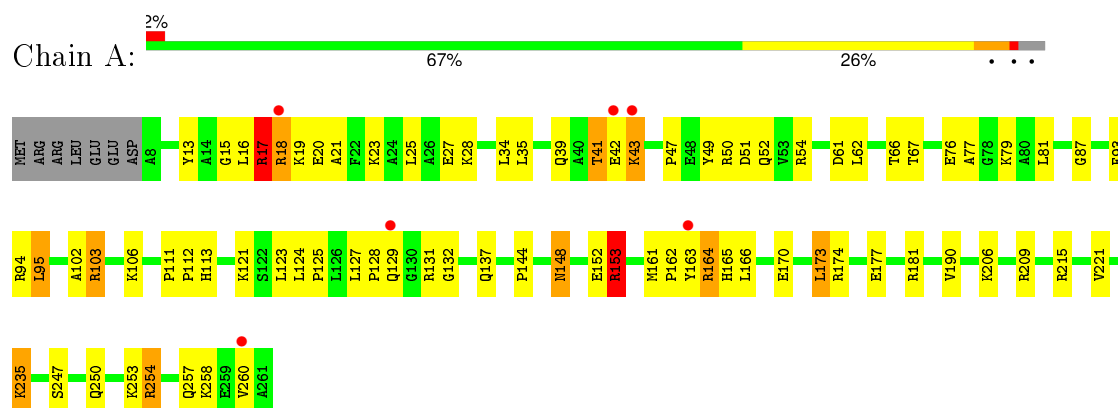
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	235	Total	O	0	0
			235	235		
2	B	228	Total	O	0	0
			228	228		

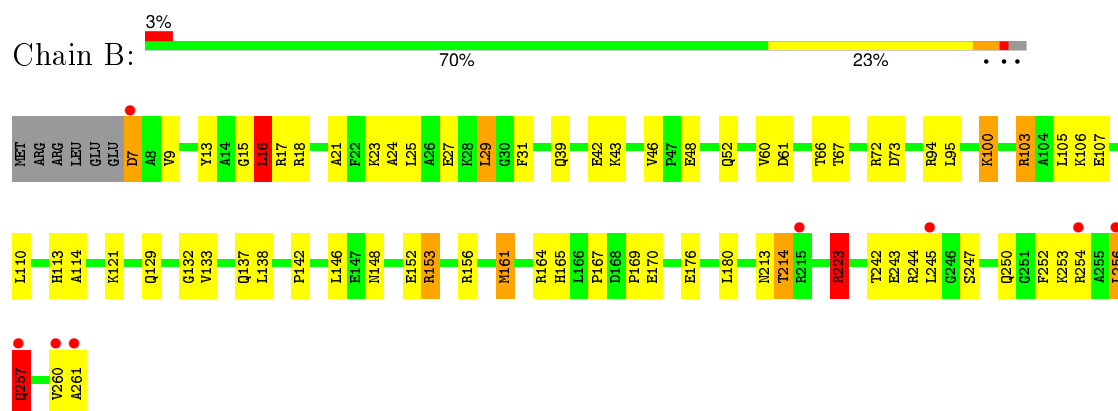
### 3 Residue-property plots

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: Uroporphyrinogen III Synthase



#### • Molecule 1: Uroporphyrinogen III Synthase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	37.90Å 84.88Å 73.95Å 90.00° 91.25° 90.00°	Depositor
Resolution (Å)	37.89 – 1.60 37.89 – 1.59	Depositor EDS
% Data completeness (in resolution range)	86.7 (37.89-1.60) 86.2 (37.89-1.59)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.58 (at 1.59Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.224 , 0.255 0.224 , 0.255	Depositor DCC
$R_{free}$ test set	2710 reflections (5.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	20.3	Xtriage
Anisotropy	0.309	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 45.5	EDS
Estimated twinning fraction	0.056 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 54690 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4357	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.38% 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	0.67	0/1979	0.88	6/2678 (0.2%)
1	B	0.74	3/1987 (0.2%)	0.92	12/2689 (0.4%)
All	All	0.70	3/3966 (0.1%)	0.90	18/5367 (0.3%)

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	A	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	252	PHE	CB-CG	-8.78	1.36	1.51
1	B	257	GLN	CA-C	6.41	1.69	1.52
1	B	257	GLN	C-O	5.87	1.34	1.23

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	164	ARG	NE-CZ-NH1	-15.07	112.77	120.30
1	A	164	ARG	NE-CZ-NH2	12.62	126.61	120.30
1	A	235	LYS	CD-CE-NZ	12.36	140.13	111.70
1	B	214	THR	N-CA-CB	-11.73	88.00	110.30
1	B	257	GLN	CB-CA-C	9.31	129.02	110.40
1	B	16	LEU	CA-CB-CG	8.99	135.98	115.30
1	A	153	ARG	NE-CZ-NH2	-8.60	116.00	120.30
1	B	256	LEU	C-N-CA	-8.50	100.44	121.70
1	B	43	LYS	N-CA-CB	-7.20	97.64	110.60
1	B	17	ARG	N-CA-C	6.53	128.63	111.00
1	B	256	LEU	N-CA-C	-6.49	93.48	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	ARG	NE-CZ-NH2	-6.13	117.23	120.30
1	A	173	LEU	CA-CB-CG	-5.84	101.86	115.30
1	B	256	LEU	CA-C-N	5.58	129.49	117.20
1	B	223	ARG	CG-CD-NE	5.44	123.23	111.80
1	B	214	THR	CB-CA-C	5.43	126.27	111.60
1	B	257	GLN	CA-C-O	5.38	131.40	120.10
1	B	252	PHE	CB-CG-CD1	-5.09	117.23	120.80

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	103	ARG	Sidechain
1	A	153	ARG	Sidechain
1	A	164	ARG	Sidechain

## 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	1943	0	2042	137	6
1	B	1951	0	2044	91	3
2	A	235	0	0	23	5
2	B	228	0	0	13	0
All	All	4357	0	4086	221	7

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 28.

All (221) 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:17:ARG:HB2	1:A:18:ARG:NE	1.42	1.33
1:B:260:VAL:CG1	1:B:261:ALA:H	1.48	1.20
1:B:244:ARG:NH2	1:B:247:SER:OG	1.79	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:LYS:NZ	2:A:494:HOH:O	1.75	1.14
1:A:177:GLU:HG2	1:A:181:ARG:HH11	1.04	1.14
1:A:173:LEU:HD21	2:A:476:HOH:O	1.46	1.10
1:B:260:VAL:HG12	1:B:261:ALA:N	1.62	1.08
1:A:17:ARG:HB2	1:A:18:ARG:CZ	1.86	1.04
1:B:94:ARG:H	1:B:113:HIS:HD2	1.06	1.03
1:B:260:VAL:HG12	1:B:261:ALA:H	0.85	1.02
1:A:18:ARG:HD2	2:A:486:HOH:O	1.61	1.01
1:B:244:ARG:CZ	1:B:247:SER:OG	2.11	0.98
1:A:173:LEU:CD2	2:A:476:HOH:O	2.04	0.97
1:A:17:ARG:CB	1:A:18:ARG:HE	1.78	0.97
1:B:129:GLN:NE2	1:B:153:ARG:HD3	1.80	0.96
1:A:17:ARG:HB3	1:A:18:ARG:HH11	1.31	0.96
1:B:29:LEU:HD23	1:B:253:LYS:HD2	1.48	0.94
1:A:177:GLU:HG2	1:A:181:ARG:NH1	1.81	0.94
1:B:244:ARG:NH1	1:B:247:SER:OG	2.01	0.93
1:A:174:ARG:NH2	2:A:267:HOH:O	2.01	0.93
1:B:25:LEU:CD2	2:B:475:HOH:O	2.15	0.93
1:A:129:GLN:HG3	1:A:153:ARG:HH21	1.34	0.91
1:A:129:GLN:CG	1:A:153:ARG:HH21	1.84	0.90
1:A:153:ARG:NH1	2:A:439:HOH:O	2.03	0.90
1:A:131:ARG:HB3	1:B:223:ARG:NH2	1.85	0.89
1:B:254:ARG:O	1:B:257:GLN:N	2.04	0.89
1:A:17:ARG:HB2	1:A:18:ARG:HE	1.11	0.88
1:A:94:ARG:H	1:A:113:HIS:HD2	1.21	0.88
1:B:25:LEU:HD22	2:B:475:HOH:O	1.74	0.88
1:A:17:ARG:CB	1:A:18:ARG:NE	2.33	0.86
1:A:19:LYS:HG2	1:A:35:LEU:HD23	1.58	0.85
1:B:107:GLU:HG2	2:B:397:HOH:O	1.75	0.84
1:A:106:LYS:HD2	1:A:112:PRO:HD2	1.59	0.84
1:B:164:ARG:NH2	2:B:336:HOH:O	2.07	0.83
1:B:29:LEU:HD23	1:B:253:LYS:CD	2.09	0.83
1:A:102:ALA:HB1	1:A:106:LYS:HZ2	1.42	0.83
1:A:206:LYS:HG2	1:A:209:ARG:NH2	1.94	0.82
1:B:261:ALA:O	2:B:398:HOH:O	1.96	0.82
1:B:129:GLN:CD	1:B:153:ARG:HD3	2.00	0.80
1:A:177:GLU:CG	1:A:181:ARG:HH11	1.92	0.80
1:B:94:ARG:H	1:B:113:HIS:CD2	1.98	0.79
1:A:94:ARG:H	1:A:113:HIS:CD2	2.01	0.79
1:B:260:VAL:CG1	1:B:261:ALA:N	2.28	0.79
1:B:213:ASN:C	1:B:214:THR:OG1	2.18	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:GLN:NE2	1:A:165:HIS:HD2	1.80	0.78
1:A:67:THR:H	1:A:137:GLN:HE22	1.31	0.77
1:A:17:ARG:CB	1:A:18:ARG:NH1	2.46	0.77
1:B:129:GLN:NE2	1:B:153:ARG:CD	2.47	0.77
1:B:29:LEU:CD2	1:B:253:LYS:HD2	2.16	0.76
1:A:17:ARG:CB	1:A:18:ARG:HH11	1.98	0.75
1:A:39:GLN:HE22	1:A:165:HIS:HD2	1.32	0.74
1:A:50:ARG:NE	1:A:54:ARG:HH12	1.85	0.74
1:A:17:ARG:CB	1:A:18:ARG:CZ	2.66	0.73
1:A:43:LYS:HD3	1:A:43:LYS:H	1.54	0.73
1:A:247:SER:HA	1:A:250:GLN:HE21	1.55	0.72
1:B:46:VAL:HG12	1:B:48:GLU:OE1	1.89	0.71
1:B:254:ARG:C	1:B:257:GLN:H	1.93	0.71
1:A:35:LEU:N	1:A:35:LEU:HD12	2.05	0.71
1:B:21:ALA:O	1:B:25:LEU:HD23	1.92	0.70
1:B:95:LEU:CD2	1:B:114:ALA:HB3	2.22	0.70
1:A:28:LYS:HD3	1:A:28:LYS:C	2.13	0.70
1:A:131:ARG:HB2	1:A:131:ARG:NH2	2.07	0.69
1:B:100:LYS:HB2	1:B:100:LYS:NZ	2.08	0.69
1:B:67:THR:H	1:B:137:GLN:HE22	1.42	0.68
1:A:28:LYS:HD3	1:A:28:LYS:O	1.92	0.68
1:A:106:LYS:HD2	1:A:112:PRO:CD	2.24	0.67
1:A:95:LEU:CD2	1:A:123:LEU:HD22	2.25	0.67
1:A:49:TYR:HA	1:A:52:GLN:HE21	1.59	0.66
1:B:73:ASP:HB3	2:B:444:HOH:O	1.94	0.66
1:A:129:GLN:HG3	1:A:153:ARG:NH2	2.11	0.66
1:A:16:LEU:HD11	1:A:39:GLN:O	1.96	0.66
1:A:13:TYR:OH	1:A:17:ARG:HG3	1.96	0.65
1:A:129:GLN:HE22	1:B:72:ARG:NH1	1.95	0.65
1:A:50:ARG:NE	1:A:54:ARG:NH1	2.44	0.65
1:A:17:ARG:HB2	1:A:18:ARG:NH1	2.12	0.64
1:A:129:GLN:CG	1:A:153:ARG:NH2	2.60	0.64
1:B:244:ARG:HH22	1:B:247:SER:HG	1.42	0.63
1:A:129:GLN:O	2:A:454:HOH:O	2.15	0.63
1:A:129:GLN:HG2	1:A:153:ARG:HH21	1.64	0.62
1:A:173:LEU:CD2	2:A:452:HOH:O	2.48	0.62
1:A:50:ARG:O	1:A:54:ARG:HG3	1.99	0.62
1:A:131:ARG:CB	1:B:223:ARG:NH2	2.61	0.62
1:A:206:LYS:HG2	1:A:209:ARG:HH21	1.63	0.62
1:A:102:ALA:HB1	1:A:106:LYS:NZ	2.15	0.62
1:B:138:LEU:HB2	1:B:161:MET:HG3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:244:ARG:HH12	1:B:247:SER:HG	1.47	0.61
1:A:131:ARG:HG2	1:B:223:ARG:HH22	1.66	0.61
1:B:167:PRO:O	1:B:169:PRO:HD3	2.01	0.61
1:B:213:ASN:C	1:B:214:THR:HG1	2.02	0.60
1:B:244:ARG:HB2	1:B:244:ARG:CZ	2.30	0.60
1:A:17:ARG:NE	1:A:190:VAL:HG21	2.16	0.60
1:A:39:GLN:NE2	1:A:165:HIS:CD2	2.68	0.59
1:A:87:GLY:HA3	1:B:254:ARG:HH12	1.67	0.59
1:A:106:LYS:HE3	2:A:326:HOH:O	2.03	0.59
1:A:50:ARG:HB3	2:A:487:HOH:O	2.02	0.59
1:A:173:LEU:HD23	2:A:452:HOH:O	2.02	0.59
1:A:95:LEU:HD22	1:A:123:LEU:HD22	1.84	0.59
1:A:17:ARG:HB3	1:A:18:ARG:NH1	2.04	0.58
1:B:24:ALA:HB3	2:B:475:HOH:O	2.04	0.58
1:B:242:THR:OG1	1:B:247:SER:HB2	2.03	0.58
1:B:142:PRO:CG	1:B:164:ARG:HH22	2.17	0.58
1:A:87:GLY:HA3	1:B:254:ARG:NH1	2.19	0.58
1:A:170:GLU:H	1:A:170:GLU:CD	2.07	0.58
1:B:39:GLN:NE2	1:B:165:HIS:ND1	2.51	0.58
1:A:124:LEU:CB	1:A:125:PRO:HD3	2.34	0.57
1:B:176:GLU:O	1:B:180:LEU:HD13	2.05	0.57
1:A:50:ARG:HG2	1:A:54:ARG:NH1	2.20	0.57
1:A:177:GLU:CG	1:A:181:ARG:NH1	2.59	0.56
1:A:177:GLU:O	1:A:181:ARG:HB2	2.05	0.56
1:A:39:GLN:HE22	1:A:165:HIS:CD2	2.19	0.56
1:A:62:LEU:HD12	1:A:93:PHE:O	2.05	0.56
1:A:18:ARG:HA	2:A:448:HOH:O	2.06	0.56
1:A:177:GLU:OE2	2:A:374:HOH:O	2.18	0.56
1:A:124:LEU:HB2	1:A:125:PRO:HD3	1.87	0.56
1:B:23:LYS:O	1:B:27:GLU:HG3	2.05	0.55
1:B:100:LYS:CB	1:B:100:LYS:NZ	2.69	0.55
1:A:260:VAL:HG23	2:A:496:HOH:O	2.05	0.55
1:A:131:ARG:HH21	1:A:131:ARG:HB2	1.70	0.55
1:A:19:LYS:CG	1:A:35:LEU:HD23	2.34	0.54
1:B:148:ASN:O	1:B:152:GLU:HG3	2.07	0.54
1:A:43:LYS:HD2	2:A:443:HOH:O	2.07	0.54
1:A:21:ALA:O	1:A:25:LEU:HD23	2.08	0.54
1:B:257:GLN:HE21	1:B:257:GLN:CA	2.21	0.53
1:B:66:THR:H	1:B:137:GLN:NE2	2.06	0.53
1:A:17:ARG:CG	1:A:18:ARG:HE	2.22	0.53
1:B:257:GLN:NE2	1:B:257:GLN:HA	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:SER:HA	1:A:250:GLN:NE2	2.22	0.52
1:B:46:VAL:CG1	1:B:48:GLU:OE1	2.56	0.52
1:A:95:LEU:HD23	1:A:123:LEU:HD22	1.92	0.52
1:B:244:ARG:CZ	1:B:244:ARG:CB	2.87	0.51
1:B:257:GLN:HE21	1:B:257:GLN:HA	1.75	0.51
1:A:13:TYR:OH	1:A:17:ARG:CG	2.58	0.51
1:A:52:GLN:NE2	2:A:331:HOH:O	2.43	0.51
1:B:42:GLU:OE2	1:B:42:GLU:N	2.41	0.51
1:B:73:ASP:CB	2:B:444:HOH:O	2.54	0.50
1:A:129:GLN:HG2	1:A:153:ARG:NH2	2.26	0.50
1:A:106:LYS:CD	1:A:112:PRO:HD2	2.35	0.50
1:A:66:THR:H	1:A:137:GLN:NE2	2.10	0.50
1:B:16:LEU:HD13	1:B:42:GLU:OE1	2.11	0.49
1:B:29:LEU:HD23	1:B:253:LYS:HD3	1.94	0.49
1:A:54:ARG:NH2	2:A:429:HOH:O	2.45	0.49
1:A:41:THR:HG23	1:A:42:GLU:H	1.76	0.49
1:B:9:VAL:HB	1:B:31:PHE:CD2	2.48	0.48
1:B:133:VAL:CG2	1:B:156:ARG:NH1	2.76	0.48
1:A:144:PRO:O	1:A:148:ASN:HB2	2.14	0.48
1:A:50:ARG:CD	1:A:54:ARG:NH1	2.76	0.48
1:B:95:LEU:HD22	1:B:114:ALA:HB3	1.92	0.48
1:B:39:GLN:OE1	1:B:165:HIS:ND1	2.46	0.48
1:B:113:HIS:HE1	2:B:487:HOH:O	1.95	0.48
1:A:42:GLU:HG3	1:A:43:LYS:N	2.29	0.48
1:B:61:ASP:OD2	1:B:132:GLY:HA3	2.14	0.48
1:A:79:LYS:HE2	2:A:470:HOH:O	2.14	0.48
1:B:129:GLN:CG	1:B:153:ARG:HD3	2.43	0.47
1:B:164:ARG:CZ	2:B:464:HOH:O	2.62	0.47
1:A:173:LEU:HD22	1:A:173:LEU:N	2.29	0.47
1:A:41:THR:HG23	1:A:42:GLU:N	2.29	0.47
1:A:253:LYS:O	1:A:257:GLN:HG3	2.16	0.46
1:B:260:VAL:HG12	1:B:261:ALA:CA	2.43	0.46
1:A:260:VAL:CG2	2:A:496:HOH:O	2.62	0.46
1:A:206:LYS:HE2	1:A:209:ARG:HD2	1.96	0.46
1:A:165:HIS:HE1	2:A:368:HOH:O	1.99	0.46
1:A:43:LYS:NZ	1:A:43:LYS:HB2	2.30	0.46
1:A:50:ARG:NE	1:A:81:LEU:CD2	2.79	0.46
1:A:50:ARG:HG2	1:A:54:ARG:CZ	2.46	0.46
1:B:152:GLU:HG3	2:B:466:HOH:O	2.15	0.46
1:B:142:PRO:HG3	1:B:164:ARG:HH22	1.81	0.46
1:A:50:ARG:HE	1:A:54:ARG:HH12	1.60	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:ALA:O	1:A:106:LYS:HD3	2.16	0.45
1:A:206:LYS:HG2	1:A:209:ARG:HH22	1.75	0.45
1:A:23:LYS:O	1:A:27:GLU:HG3	2.17	0.45
1:B:170:GLU:H	1:B:170:GLU:CD	2.20	0.45
1:B:13:TYR:CZ	1:B:15:GLY:HA3	2.52	0.45
1:B:103:ARG:HE	1:B:103:ARG:HB2	1.47	0.45
1:B:142:PRO:CG	1:B:164:ARG:NH2	2.80	0.45
1:B:60:VAL:HG12	1:B:133:VAL:CG1	2.47	0.45
1:A:128:PRO:O	1:A:153:ARG:NE	2.42	0.45
1:A:93:PHE:HE2	1:A:95:LEU:HD11	1.82	0.45
1:A:206:LYS:HE2	1:A:209:ARG:HH21	1.80	0.44
1:A:93:PHE:CE2	1:A:95:LEU:HD12	2.52	0.44
1:A:76:GLU:HG2	1:A:77:ALA:N	2.31	0.44
1:A:41:THR:CG2	1:A:42:GLU:N	2.81	0.44
1:A:51:ASP:HA	1:A:54:ARG:HG3	2.00	0.44
1:A:61:ASP:OD2	1:A:132:GLY:HA3	2.17	0.44
1:B:52:GLN:NE2	2:B:340:HOH:O	2.50	0.44
1:A:34:LEU:C	1:A:35:LEU:HD12	2.38	0.43
1:B:39:GLN:CD	1:B:165:HIS:ND1	2.72	0.43
1:B:9:VAL:HB	1:B:31:PHE:HD2	1.83	0.43
1:B:25:LEU:O	1:B:29:LEU:HD12	2.17	0.43
1:A:106:LYS:NZ	2:A:286:HOH:O	2.47	0.43
1:A:15:GLY:HA3	1:A:17:ARG:HD3	2.00	0.43
1:A:190:VAL:HA	1:A:221:VAL:HB	1.99	0.43
1:B:100:LYS:HB2	1:B:100:LYS:HZ2	1.83	0.43
1:B:105:LEU:HB3	1:B:110:LEU:O	2.19	0.43
1:A:131:ARG:CG	1:B:223:ARG:HH22	2.31	0.43
1:A:17:ARG:CG	1:A:18:ARG:H	2.32	0.42
1:A:127:LEU:O	1:A:153:ARG:CZ	2.67	0.42
1:B:60:VAL:HG12	1:B:133:VAL:HG11	2.01	0.42
1:A:173:LEU:CD1	2:A:476:HOH:O	2.67	0.42
1:A:50:ARG:HE	1:A:54:ARG:NH1	2.16	0.42
1:A:62:LEU:HD13	1:A:93:PHE:HD2	1.83	0.42
1:A:161:MET:HG2	1:A:163:TYR:O	2.19	0.42
1:B:142:PRO:HG3	1:B:164:ARG:NH2	2.35	0.42
1:A:93:PHE:CE2	1:A:95:LEU:CD1	3.03	0.42
1:A:20:GLU:OE2	1:A:20:GLU:HA	2.20	0.41
1:A:127:LEU:HA	1:A:128:PRO:HD3	1.81	0.41
1:A:106:LYS:CE	2:A:326:HOH:O	2.65	0.41
1:A:28:LYS:C	1:A:28:LYS:CD	2.84	0.41
1:B:133:VAL:HG22	1:B:156:ARG:NH1	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:GLU:CG	1:A:77:ALA:N	2.83	0.41
1:B:106:LYS:CB	1:B:106:LYS:NZ	2.84	0.41
1:A:93:PHE:HE2	1:A:95:LEU:CD1	2.33	0.41
1:A:161:MET:HA	1:A:162:PRO:HD2	1.95	0.41
1:B:242:THR:C	1:B:243:GLU:HG3	2.41	0.41
1:B:244:ARG:NH1	1:B:247:SER:HG	2.08	0.41
1:A:43:LYS:N	1:A:43:LYS:HD3	2.29	0.41
1:A:235:LYS:HD2	1:A:235:LYS:HA	1.56	0.41
1:B:7:ASP:OD2	1:B:7:ASP:O	2.39	0.41
1:B:121:LYS:HE3	2:B:481:HOH:O	2.21	0.41
1:A:106:LYS:HG3	1:A:111:PRO:HB3	2.02	0.40
1:A:166:LEU:HA	1:A:166:LEU:HD23	1.85	0.40
1:B:18:ARG:NH2	1:B:245:LEU:HD12	2.36	0.40

All (7) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:GLU:OE1	1:B:254:ARG:CZ[2_645]	1.72	0.48
1:B:106:LYS:NZ	2:A:282:HOH:O[1_455]	1.82	0.38
1:A:258:LYS:NZ	2:A:464:HOH:O[1_655]	1.91	0.29
1:A:28:LYS:CE	2:A:457:HOH:O[1_655]	1.95	0.25
1:A:28:LYS:NZ	2:A:457:HOH:O[1_655]	1.96	0.24
1:A:20:GLU:OE1	1:B:254:ARG:NH1[2_645]	1.98	0.22
1:A:254:ARG:NH2	2:A:475:HOH:O[1_655]	1.99	0.21

## 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	252/261 (97%)	247 (98%)	4 (2%)	1 (0%)	39 17

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	253/261 (97%)	248 (98%)	5 (2%)	0	100	100
All	All	505/522 (97%)	495 (98%)	9 (2%)	1 (0%)	52	28

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	17	ARG

### 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	192/199 (96%)	182 (95%)	10 (5%)	29	7
1	B	193/199 (97%)	181 (94%)	12 (6%)	23	5
All	All	385/398 (97%)	363 (94%)	22 (6%)	25	6

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ARG
1	A	18	ARG
1	A	41	THR
1	A	43	LYS
1	A	47	PRO
1	A	95	LEU
1	A	148	ASN
1	A	152	GLU
1	A	215	ARG
1	A	254	ARG
1	B	7	ASP
1	B	16	LEU
1	B	29	LEU
1	B	100	LYS
1	B	103	ARG

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Mol	Chain	Res	Type
1	B	146	LEU
1	B	153	ARG
1	B	161	MET
1	B	223	ARG
1	B	250	GLN
1	B	256	LEU
1	B	257	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	52	GLN
1	A	113	HIS
1	A	129	GLN
1	A	137	GLN
1	A	148	ASN
1	A	165	HIS
1	A	250	GLN
1	B	39	GLN
1	B	52	GLN
1	B	113	HIS
1	B	137	GLN
1	B	250	GLN
1	B	257	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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

There are no ligands in this entry.

## 5.7 Other polymers

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(Å <sup>2</sup> )	Q<0.9
1	A	254/261 (97%)	-0.17	6 (2%) 62 60	13, 25, 46, 59	0
1	B	255/261 (97%)	-0.15	8 (3%) 52 50	15, 26, 48, 73	0
All	All	509/522 (97%)	-0.16	14 (2%) 56 54	13, 26, 47, 73	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	129	GLN	5.0
1	B	260	VAL	4.3
1	A	260	VAL	4.0
1	B	261	ALA	3.8
1	A	43	LYS	3.0
1	B	257	GLN	2.9
1	B	7	ASP	2.7
1	B	254	ARG	2.3
1	B	256	LEU	2.2
1	B	215	ARG	2.2
1	A	18	ARG	2.1
1	A	163	TYR	2.1
1	A	42	GLU	2.0
1	B	245	LEU	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 6.3 Carbohydrates [i](#)

There are no 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.