Full wwPDB X-ray Structure Validation Report (i)

Aug 15, 2023 – 09:57 PM EDT

PDB ID	:	1XRT
Title	:	The Crystal Structure of a Novel, Latent Dihydroorotase from Aquifex Aeolicus at 1.7 A Resolution
Authors	:	Martin, P.D.; Purcarea, C.; Zhang, P.; Vaishnav, A.; Sadecki, S.; Guy-Evans, H.I.; Evans, D.R.; Edwards, B.F.
Deposited on	:	2004-10-15
Resolution	:	1.61 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

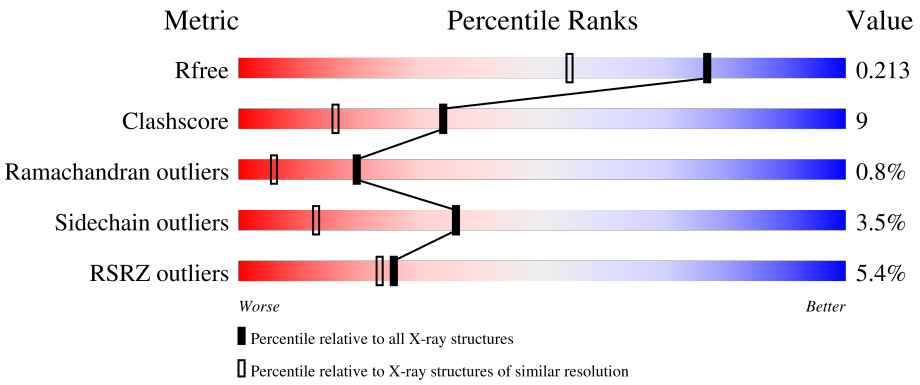
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.61 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (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 of 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 <=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	467	5% 67% 10% • 21%
1	B	467	4% 66% 12% • 21%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6940 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 Dihydroorotase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	369	2849	1815	482	537	15	0	0	0
1	B	369	2850	1818	483	534	15	0	0	0

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-44	MET	-	cloning artifact	UNP O66990
A	-43	ARG	-	cloning artifact	UNP O66990
A	-42	PHE	-	cloning artifact	UNP O66990
A	-41	TRP	-	cloning artifact	UNP O66990
A	-40	GLN	-	cloning artifact	UNP O66990
A	-39	TYR	-	cloning artifact	UNP O66990
A	-38	ILE	-	cloning artifact	UNP O66990
A	-37	ASN	-	cloning artifact	UNP O66990
A	-36	GLY	-	cloning artifact	UNP O66990
A	-35	VAL	-	cloning artifact	UNP O66990
A	-34	ASP	-	cloning artifact	UNP O66990
A	-33	MET	-	cloning artifact	UNP O66990
A	-32	ARG	-	cloning artifact	UNP O66990
A	-31	GLY	-	cloning artifact	UNP O66990
A	-30	SER	-	cloning artifact	UNP O66990
A	-29	HIS	-	cloning artifact	UNP O66990
A	-28	HIS	-	cloning artifact	UNP O66990
A	-27	HIS	-	cloning artifact	UNP O66990
A	-26	HIS	-	cloning artifact	UNP O66990
A	-25	HIS	-	cloning artifact	UNP O66990
A	-24	HIS	-	cloning artifact	UNP O66990
A	-23	GLY	-	cloning artifact	UNP O66990
A	-22	MET	-	cloning artifact	UNP O66990
A	-21	ALA	-	cloning artifact	UNP O66990
A	-20	SER	-	cloning artifact	UNP O66990

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	cloning artifact	UNP O66990
A	-18	THR	-	cloning artifact	UNP O66990
A	-17	GLY	-	cloning artifact	UNP O66990
A	-16	GLY	-	cloning artifact	UNP O66990
A	-15	GLN	-	cloning artifact	UNP O66990
A	-14	GLN	-	cloning artifact	UNP O66990
A	-13	MET	-	cloning artifact	UNP O66990
A	-12	GLY	-	cloning artifact	UNP O66990
A	-11	ARG	-	cloning artifact	UNP O66990
A	-10	ASP	-	cloning artifact	UNP O66990
A	-9	LEU	-	cloning artifact	UNP O66990
A	-8	TYR	-	cloning artifact	UNP O66990
A	-7	ASP	-	cloning artifact	UNP O66990
A	-6	ASP	-	cloning artifact	UNP O66990
A	-5	ASP	-	cloning artifact	UNP O66990
A	-4	ASP	-	cloning artifact	UNP O66990
A	-3	LYS	-	cloning artifact	UNP O66990
A	-2	ASP	-	cloning artifact	UNP O66990
A	-1	ARG	-	cloning artifact	UNP O66990
A	0	TRP	-	cloning artifact	UNP O66990
B	-44	MET	-	cloning artifact	UNP O66990
B	-43	ARG	-	cloning artifact	UNP O66990
B	-42	PHE	-	cloning artifact	UNP O66990
B	-41	TRP	-	cloning artifact	UNP O66990
B	-40	GLN	-	cloning artifact	UNP O66990
B	-39	TYR	-	cloning artifact	UNP O66990
B	-38	ILE	-	cloning artifact	UNP O66990
B	-37	ASN	-	cloning artifact	UNP O66990
B	-36	GLY	-	cloning artifact	UNP O66990
B	-35	VAL	-	cloning artifact	UNP O66990
B	-34	ASP	-	cloning artifact	UNP O66990
B	-33	MET	-	cloning artifact	UNP O66990
B	-32	ARG	-	cloning artifact	UNP O66990
B	-31	GLY	-	cloning artifact	UNP O66990
B	-30	SER	-	cloning artifact	UNP O66990
B	-29	HIS	-	cloning artifact	UNP O66990
B	-28	HIS	-	cloning artifact	UNP O66990
B	-27	HIS	-	cloning artifact	UNP O66990
B	-26	HIS	-	cloning artifact	UNP O66990
B	-25	HIS	-	cloning artifact	UNP O66990
B	-24	HIS	-	cloning artifact	UNP O66990
B	-23	GLY	-	cloning artifact	UNP O66990

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	MET	-	cloning artifact	UNP O66990
B	-21	ALA	-	cloning artifact	UNP O66990
B	-20	SER	-	cloning artifact	UNP O66990
B	-19	MET	-	cloning artifact	UNP O66990
B	-18	THR	-	cloning artifact	UNP O66990
B	-17	GLY	-	cloning artifact	UNP O66990
B	-16	GLY	-	cloning artifact	UNP O66990
B	-15	GLN	-	cloning artifact	UNP O66990
B	-14	GLN	-	cloning artifact	UNP O66990
B	-13	MET	-	cloning artifact	UNP O66990
B	-12	GLY	-	cloning artifact	UNP O66990
B	-11	ARG	-	cloning artifact	UNP O66990
B	-10	ASP	-	cloning artifact	UNP O66990
B	-9	LEU	-	cloning artifact	UNP O66990
B	-8	TYR	-	cloning artifact	UNP O66990
B	-7	ASP	-	cloning artifact	UNP O66990
B	-6	ASP	-	cloning artifact	UNP O66990
B	-5	ASP	-	cloning artifact	UNP O66990
B	-4	ASP	-	cloning artifact	UNP O66990
B	-3	LYS	-	cloning artifact	UNP O66990
B	-2	ASP	-	cloning artifact	UNP O66990
B	-1	ARG	-	cloning artifact	UNP O66990
B	0	TRP	-	cloning artifact	UNP O66990

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	A	1	1	1	0	0
			Total	Zn		
2	B	1	1	1	0	0

- Molecule 3 is water.

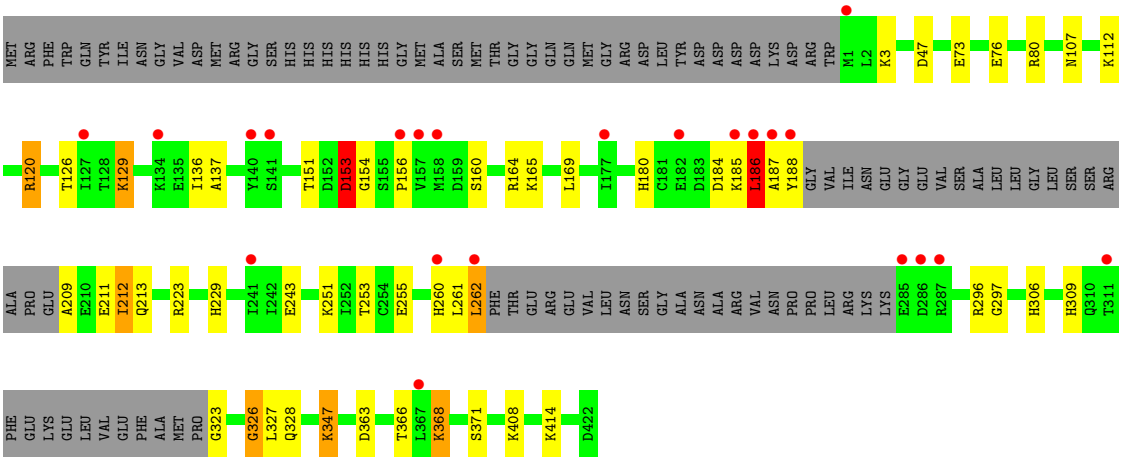
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	625	625	625	0	0
			Total	O		
3	B	614	614	614	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

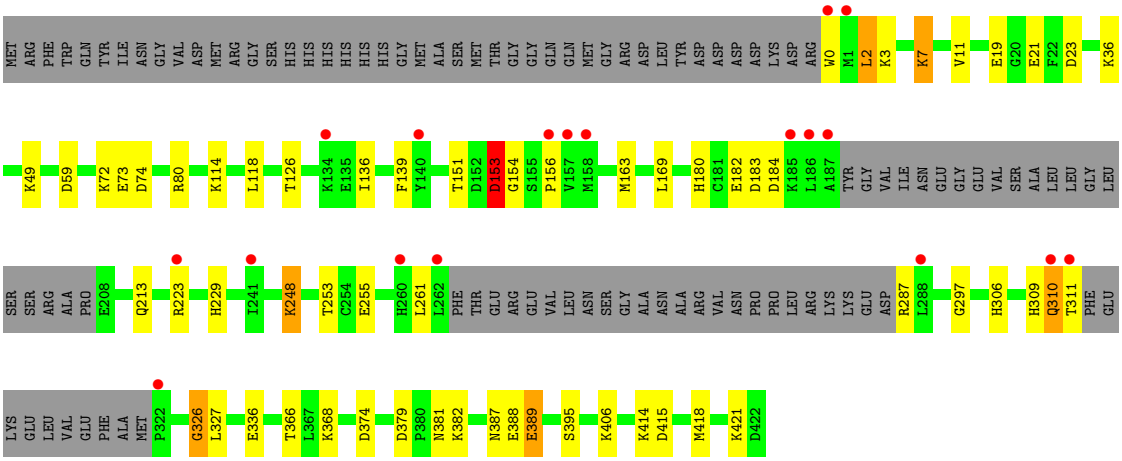
- Molecule 1: Dihydroorotase

Chain A:



- Molecule 1: Dihydroorotase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	174.20Å 91.23Å 61.18Å 90.00° 93.15° 90.00°	Depositor
Resolution (Å)	20.00 – 1.61 19.95 – 1.61	Depositor EDS
% Data completeness (in resolution range)	77.6 (20.00-1.61) 77.6 (19.95-1.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$< I/\sigma(I) >$ [1]	2.40 (at 1.61Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.165 , 0.208 0.172 , 0.213	Depositor DCC
R_{free} test set	4873 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	26.0	Xtriage
Anisotropy	0.373	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 44.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	6940	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.01% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.87	0/2893	0.91	3/3904 (0.1%)
1	B	0.90	1/2896 (0.0%)	0.95	7/3909 (0.2%)
All	All	0.88	1/5789 (0.0%)	0.93	10/7813 (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	A	0	3
1	B	0	1
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	163	MET	SD-CE	-5.69	1.46	1.77

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	415	ASP	CB-CG-OD2	6.87	124.49	118.30
1	B	74	ASP	CB-CG-OD1	6.73	124.36	118.30
1	A	153	ASP	CB-CG-OD1	6.53	124.18	118.30
1	B	2	LEU	CB-CG-CD2	6.09	121.36	111.00
1	B	59	ASP	CB-CG-OD2	5.91	123.61	118.30
1	B	374	ASP	CB-CG-OD2	5.74	123.46	118.30
1	A	47	ASP	CB-CG-OD2	5.42	123.18	118.30
1	B	183	ASP	CB-CG-OD2	5.37	123.14	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	363	ASP	CB-CG-OD2	5.36	123.12	118.30
1	B	379	ASP	CB-CG-OD2	5.10	122.89	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	153	ASP	Peptide
1	A	185	LYS	Peptide
1	A	186	LEU	Peptide
1	B	153	ASP	Peptide

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2849	0	2945	57	0
1	B	2850	0	2950	42	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	625	0	0	28	1
3	B	614	0	0	12	1
All	All	6940	0	5895	99	1

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 9.

All (99) 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:371:SER:HB2	3:A:1872:HOH:O	1.09	1.27
1:A:366:THR:HG22	3:A:1846:HOH:O	1.42	1.20
1:A:186:LEU:O	1:A:309:HIS:HE1	1.26	1.15
1:A:186:LEU:O	1:A:309:HIS:CE1	2.01	1.12
1:B:156:PRO:HB3	1:B:213:GLN:HE22	1.18	1.08

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:LEU:C	3:A:1857:HOH:O	1.95	1.03
1:B:366:THR:HG22	3:B:1475:HOH:O	1.63	0.99
1:A:209:ALA:N	3:A:1859:HOH:O	2.04	0.91
1:A:323:GLY:N	3:A:1867:HOH:O	2.08	0.87
1:B:156:PRO:HB3	1:B:213:GLN:NE2	1.92	0.84
1:B:366:THR:HG23	1:B:368:LYS:H	1.44	0.82
1:A:251:LYS:HD3	3:A:1589:HOH:O	1.80	0.82
1:B:136:ILE:HD11	1:B:169:LEU:HD12	1.61	0.81
1:B:387:ASN:OD1	1:B:389:GLU:HG2	1.86	0.75
1:A:76:GLU:OE1	1:A:80:ARG:NH2	2.20	0.74
1:A:129:LYS:NZ	3:A:1874:HOH:O	2.20	0.74
1:B:156:PRO:CB	1:B:213:GLN:HE22	1.97	0.73
1:B:3:LYS:HD3	3:B:1861:HOH:O	1.88	0.73
1:A:212:ILE:HB	3:A:1866:HOH:O	1.88	0.73
1:A:120:ARG:HD2	3:A:1478:HOH:O	1.90	0.71
1:B:0:TRP:N	3:B:1874:HOH:O	2.21	0.70
1:B:72:LYS:HG2	1:B:310:GLN:HG2	1.73	0.70
1:A:107:ASN:HB3	3:A:1818:HOH:O	1.89	0.70
1:A:229:HIS:HE1	1:A:255:GLU:OE1	1.75	0.69
1:A:186:LEU:HD12	3:A:1855:HOH:O	1.93	0.68
1:A:309:HIS:ND1	3:A:1858:HOH:O	2.27	0.67
1:A:297:GLY:HA2	3:A:1555:HOH:O	1.94	0.67
1:B:309:HIS:HE1	3:B:1582:HOH:O	1.79	0.66
1:A:136:ILE:HD11	1:A:169:LEU:HD12	1.77	0.65
1:B:229:HIS:HE1	1:B:255:GLU:OE1	1.79	0.65
1:A:156:PRO:HB3	1:A:213:GLN:HE22	1.63	0.64
1:A:211:GLU:CD	3:A:1861:HOH:O	2.37	0.63
1:B:229:HIS:HD2	1:B:253:THR:OG1	1.82	0.62
1:A:328:GLN:HG2	3:A:1876:HOH:O	2.00	0.62
1:A:164:ARG:NH2	1:A:223:ARG:HG2	2.17	0.59
1:B:310:GLN:NE2	1:B:395:SER:OG	2.36	0.59
1:A:306:HIS:HD2	1:A:326:GLY:H	1.50	0.59
1:B:21:GLU:O	1:B:36:LYS:HD3	2.03	0.59
1:A:306:HIS:CD2	1:A:326:GLY:H	2.20	0.58
1:A:76:GLU:OE1	1:A:80:ARG:CZ	2.51	0.58
1:B:421:LYS:NZ	3:B:1622:HOH:O	2.35	0.58
1:A:154:GLY:CA	1:A:184:ASP:OD1	2.52	0.58
1:B:306:HIS:HD2	1:B:327:LEU:H	1.50	0.57
1:A:366:THR:HG23	1:A:368:LYS:H	1.69	0.57
1:A:408:LYS:HE2	3:A:1750:HOH:O	2.04	0.57
1:B:297:GLY:HA2	3:B:1639:HOH:O	2.05	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:HIS:CD2	3:A:1571:HOH:O	2.59	0.56
1:B:306:HIS:HD2	1:B:326:GLY:H	1.53	0.56
1:B:388:GLU:HG2	1:B:389:GLU:OE1	2.06	0.55
1:B:306:HIS:CD2	1:B:326:GLY:H	2.24	0.55
1:A:306:HIS:HD2	1:A:327:LEU:H	1.54	0.54
1:B:49:LYS:HE2	3:B:1788:HOH:O	2.05	0.54
1:A:243:GLU:HG3	3:A:1550:HOH:O	2.06	0.54
1:A:414:LYS:NZ	3:A:1865:HOH:O	1.96	0.54
1:B:7:LYS:NZ	1:B:23:ASP:OD2	2.38	0.54
1:A:186:LEU:HD22	1:A:186:LEU:C	2.28	0.53
1:B:118:LEU:O	1:B:414:LYS:HD3	2.09	0.52
1:B:154:GLY:CA	1:B:184:ASP:OD1	2.58	0.52
1:A:154:GLY:HA3	1:A:184:ASP:OD1	2.10	0.52
1:A:186:LEU:HD13	1:A:309:HIS:ND1	2.25	0.52
1:B:80:ARG:NH1	3:B:1864:HOH:O	2.43	0.51
1:B:153:ASP:OD1	1:B:154:GLY:N	2.44	0.50
1:A:309:HIS:HB3	3:A:1855:HOH:O	2.11	0.50
1:A:243:GLU:OE2	1:A:296:ARG:NH2	2.45	0.49
1:B:154:GLY:HA2	1:B:184:ASP:OD1	2.13	0.49
1:A:165:LYS:NZ	1:A:169:LEU:HG	2.26	0.49
1:A:229:HIS:HD2	1:A:253:THR:OG1	1.95	0.49
1:A:186:LEU:HD21	3:A:1595:HOH:O	2.12	0.49
1:A:209:ALA:CA	3:A:1859:HOH:O	2.57	0.49
1:B:154:GLY:CA	1:B:184:ASP:CG	2.81	0.49
1:A:73:GLU:OE2	1:A:306:HIS:HE1	1.95	0.49
1:A:186:LEU:CD1	3:A:1855:HOH:O	2.57	0.48
1:A:154:GLY:HA2	1:A:184:ASP:OD1	2.14	0.48
1:B:73:GLU:OE2	1:B:306:HIS:HE1	1.96	0.48
1:B:248:LYS:HG2	3:B:1571:HOH:O	2.14	0.48
1:A:153:ASP:OD1	1:A:154:GLY:N	2.47	0.47
1:B:139:PHE:HB2	1:B:169:LEU:HD13	1.97	0.46
1:A:112:LYS:HE3	3:A:1458:HOH:O	2.14	0.46
1:A:186:LEU:C	1:A:186:LEU:CD2	2.83	0.46
1:A:154:GLY:CA	1:A:184:ASP:CG	2.84	0.46
1:B:366:THR:OG1	1:B:368:LYS:HG3	2.16	0.46
1:A:129:LYS:HD2	1:A:137:ALA:HA	2.00	0.44
1:A:126:THR:HA	1:A:151:THR:O	2.17	0.44
1:A:136:ILE:HG22	3:A:1737:HOH:O	2.16	0.44
1:A:347:LYS:HE3	3:A:1790:HOH:O	2.18	0.44
1:B:114:LYS:HB3	1:B:114:LYS:NZ	2.32	0.43
1:A:261:LEU:O	1:A:262:LEU:HB2	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:ALA:O	1:A:212:ILE:HG22	2.18	0.43
1:B:0:TRP:CA	3:B:1874:HOH:O	2.65	0.43
1:B:381:ASN:HD22	1:B:406:LYS:HE2	1.83	0.42
1:A:187:ALA:HB3	1:A:188:TYR:CD1	2.55	0.42
1:B:154:GLY:HA3	1:B:184:ASP:OD1	2.19	0.42
1:B:49:LYS:CE	3:B:1788:HOH:O	2.65	0.41
1:A:251:LYS:HD2	3:A:1774:HOH:O	2.21	0.41
1:A:328:GLN:CG	3:A:1876:HOH:O	2.65	0.41
1:B:126:THR:HA	1:B:151:THR:O	2.21	0.41
1:B:306:HIS:CD2	1:B:327:LEU:H	2.34	0.41
1:B:7:LYS:NZ	3:B:1605:HOH:O	2.37	0.40
1:B:11:VAL:O	1:B:19:GLU:HA	2.22	0.40

All (1) 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)
3:A:1493:HOH:O	3:B:1889:HOH:O[4_545]	2.14	0.06

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	361/467 (77%)	342 (95%)	16 (4%)	3 (1%)	19 6
1	B	361/467 (77%)	343 (95%)	15 (4%)	3 (1%)	19 6
All	All	722/934 (77%)	685 (95%)	31 (4%)	6 (1%)	19 6

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	153	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	153	ASP
1	A	180	HIS
1	B	326	GLY
1	A	326	GLY
1	B	180	HIS

5.3.2 Protein sidechains (i)

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	318/400 (80%)	309 (97%)	9 (3%)	43 18
1	B	318/400 (80%)	305 (96%)	13 (4%)	30 9
All	All	636/800 (80%)	614 (96%)	22 (4%)	36 13

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

Mol	Chain	Res	Type
1	A	3	LYS
1	A	120	ARG
1	A	129	LYS
1	A	160	SER
1	A	186	LEU
1	A	212	ILE
1	A	262	LEU
1	A	347	LYS
1	A	368	LYS
1	B	2	LEU
1	B	7	LYS
1	B	182	GLU
1	B	223	ARG
1	B	248	LYS
1	B	261	LEU
1	B	287	ARG
1	B	310	GLN
1	B	311	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	336	GLU
1	B	382	LYS
1	B	389	GLU
1	B	418	MET

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

Mol	Chain	Res	Type
1	A	16	GLN
1	A	213	GLN
1	A	229	HIS
1	A	306	HIS
1	A	309	HIS
1	A	381	ASN
1	B	16	GLN
1	B	180	HIS
1	B	213	GLN
1	B	229	HIS
1	B	231	GLN
1	B	306	HIS
1	B	310	GLN
1	B	381	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	369/467 (79%)	0.30	22 (5%)	21	19	12, 18, 35, 57	0
1	B	369/467 (79%)	0.35	18 (4%)	29	27	12, 18, 34, 51	0
All	All	738/934 (79%)	0.32	40 (5%)	25	23	12, 18, 35, 57	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	186	LEU	11.1
1	A	186	LEU	10.9
1	B	311	THR	8.1
1	A	285	GLU	6.0
1	A	187	ALA	5.1
1	B	322	PRO	5.0
1	B	187	ALA	4.8
1	B	262	LEU	4.4
1	A	185	LYS	4.4
1	B	157	VAL	4.3
1	A	188	TYR	4.0
1	A	182	GLU	3.9
1	B	310	GLN	3.8
1	A	157	VAL	3.8
1	A	140	TYR	3.8
1	A	262	LEU	3.6
1	B	158	MET	3.5
1	B	156	PRO	3.4
1	A	1	MET	3.4
1	A	134	LYS	3.4
1	A	156	PRO	3.3
1	B	185	LYS	3.2
1	A	177	ILE	2.9
1	B	1	MET	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	311	THR	2.8
1	B	134	LYS	2.7
1	A	260	HIS	2.7
1	A	286	ASP	2.6
1	B	0	TRP	2.6
1	B	140	TYR	2.5
1	A	158	MET	2.5
1	A	127	ILE	2.5
1	B	260	HIS	2.5
1	A	367	LEU	2.4
1	B	288	LEU	2.3
1	B	241	ILE	2.3
1	A	141	SER	2.2
1	A	287	ARG	2.1
1	A	241	ILE	2.1
1	B	223	ARG	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 monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ZN	A	1423	1/1	1.00	0.08	34,34,34,34	0
2	ZN	B	1424	1/1	1.00	0.08	33,33,33,33	0

6.5 Other polymers (i)

There are no such residues in this entry.