



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 12:22 pm BST

PDB ID : 4WL3
Title : Crystal structure determination of Bile Salt Hydrolase from *Enterococcus faecalis*
Authors : Ramasamy, S.; Chand, D.; Suresh, C.G.
Deposited on : 2014-10-06
Resolution : 2.01 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

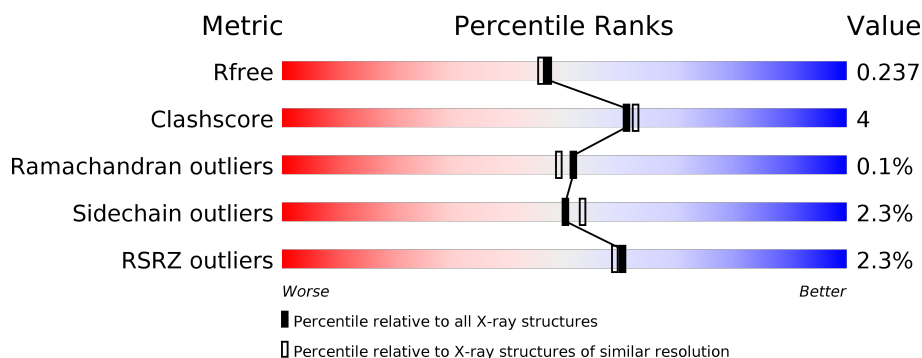
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	331	<div> <div>3%</div> <div> <div></div> <div>85%</div> <div>11%</div> <div>..</div> </div> </div>
1	B	331	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>9%</div> <div>..</div> </div> </div>
1	C	331	<div> <div>2%</div> <div> <div></div> <div>89%</div> <div>8%</div> <div>..</div> </div> </div>
1	D	331	<div> <div>2%</div> <div> <div></div> <div>88%</div> <div>8%</div> <div>...</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10507 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 Bile salt hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	0	0
			2597	1658	420	510	9			
1	B	323	Total	C	N	O	S	0	0	0
			2597	1658	420	510	9			
1	C	323	Total	C	N	O	S	0	0	0
			2597	1658	420	510	9			
1	D	323	Total	C	N	O	S	0	0	0
			2597	1658	420	510	9			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	325	LEU	-	expression tag	UNP C7CXJ5
A	326	GLU	-	expression tag	UNP C7CXJ5
A	327	HIS	-	expression tag	UNP C7CXJ5
A	328	HIS	-	expression tag	UNP C7CXJ5
A	329	HIS	-	expression tag	UNP C7CXJ5
A	330	HIS	-	expression tag	UNP C7CXJ5
A	331	HIS	-	expression tag	UNP C7CXJ5
A	332	HIS	-	expression tag	UNP C7CXJ5
B	325	LEU	-	expression tag	UNP C7CXJ5
B	326	GLU	-	expression tag	UNP C7CXJ5
B	327	HIS	-	expression tag	UNP C7CXJ5
B	328	HIS	-	expression tag	UNP C7CXJ5
B	329	HIS	-	expression tag	UNP C7CXJ5
B	330	HIS	-	expression tag	UNP C7CXJ5
B	331	HIS	-	expression tag	UNP C7CXJ5
B	332	HIS	-	expression tag	UNP C7CXJ5
C	325	LEU	-	expression tag	UNP C7CXJ5
C	326	GLU	-	expression tag	UNP C7CXJ5
C	327	HIS	-	expression tag	UNP C7CXJ5
C	328	HIS	-	expression tag	UNP C7CXJ5
C	329	HIS	-	expression tag	UNP C7CXJ5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	330	HIS	-	expression tag	UNP C7CXJ5
C	331	HIS	-	expression tag	UNP C7CXJ5
C	332	HIS	-	expression tag	UNP C7CXJ5
D	325	LEU	-	expression tag	UNP C7CXJ5
D	326	GLU	-	expression tag	UNP C7CXJ5
D	327	HIS	-	expression tag	UNP C7CXJ5
D	328	HIS	-	expression tag	UNP C7CXJ5
D	329	HIS	-	expression tag	UNP C7CXJ5
D	330	HIS	-	expression tag	UNP C7CXJ5
D	331	HIS	-	expression tag	UNP C7CXJ5
D	332	HIS	-	expression tag	UNP C7CXJ5

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	34	Total O 34 34	0	0
2	B	30	Total O 30 30	0	0
2	C	32	Total O 32 32	0	0
2	D	23	Total O 23 23	0	0

- Molecule 1: Bile salt hydrolase



K306 E316 I320 N324 LEU GLU HIS HIS HIS HIS HIS HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.20 Å 131.62 Å 86.72 Å 90.00° 94.48° 90.00°	Depositor
Resolution (Å)	41.07 – 2.01 41.07 – 2.01	Depositor EDS
% Data completeness (in resolution range)	95.3 (41.07-2.01) 95.4 (41.07-2.01)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.64 (at 2.01 Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.205 , 0.235 0.212 , 0.237	Depositor DCC
R_{free} test set	4645 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	16.2	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10507	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% 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	0.97	0/2656	0.92	4/3601 (0.1%)
1	B	0.93	1/2656 (0.0%)	0.95	5/3601 (0.1%)
1	C	0.94	1/2656 (0.0%)	1.04	4/3601 (0.1%)
1	D	0.91	1/2656 (0.0%)	1.04	8/3601 (0.2%)
All	All	0.94	3/10624 (0.0%)	0.99	21/14404 (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	1
1	B	0	1
1	D	0	1
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	57	GLY	N-CA	7.61	1.57	1.46
1	D	223	ARG	CD-NE	-6.06	1.36	1.46
1	C	223	ARG	CD-NE	-5.51	1.37	1.46

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	223	ARG	NE-CZ-NH2	-22.93	108.83	120.30
1	D	223	ARG	NE-CZ-NH2	-22.37	109.12	120.30
1	C	223	ARG	NE-CZ-NH1	21.61	131.10	120.30
1	D	223	ARG	NE-CZ-NH1	18.83	129.72	120.30
1	D	223	ARG	CD-NE-CZ	9.30	136.62	123.60
1	A	223	ARG	NE-CZ-NH2	-8.91	115.85	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	56	ALA	C-N-CA	-8.76	103.89	122.30
1	D	306	LYS	CD-CE-NZ	7.44	128.82	111.70
1	C	223	ARG	CD-NE-CZ	7.12	133.56	123.60
1	B	223	ARG	NE-CZ-NH2	-7.10	116.75	120.30
1	D	288	THR	N-CA-CB	-6.97	97.06	110.30
1	C	324	ASN	N-CA-CB	-6.30	99.27	110.60
1	A	147	GLN	N-CA-CB	-6.23	99.39	110.60
1	B	226	LYS	CD-CE-NZ	6.22	126.00	111.70
1	A	135	LEU	CB-CG-CD1	5.96	121.14	111.00
1	D	187	LEU	CA-CB-CG	5.58	128.13	115.30
1	B	226	LYS	CG-CD-CE	5.58	128.63	111.90
1	B	260	LEU	CB-CG-CD1	-5.31	101.97	111.00
1	D	65	TYR	CB-CG-CD1	5.24	124.14	121.00
1	A	307	ASP	CB-CA-C	-5.22	99.97	110.40
1	D	40	ARG	NE-CZ-NH1	5.04	122.82	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	127	ILE	Peptide
1	B	57	GLY	Peptide
1	D	223	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	2597	0	2504	33	1
1	B	2597	0	2504	30	0
1	C	2597	0	2504	19	1
1	D	2597	0	2504	29	0
2	A	34	0	0	0	1
2	B	30	0	0	0	0
2	C	32	0	0	0	1
2	D	23	0	0	0	0
All	All	10507	0	10016	91	2

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 4.

All (91) 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:218:LEU:O	1:A:223:ARG:NH1	1.66	1.26
1:A:207:ARG:NH1	1:B:21:GLU:OE2	1.98	0.96
1:A:75:MET:HE1	1:A:105:VAL:HG11	1.56	0.85
1:D:75:MET:HE1	1:D:105:VAL:HG11	1.56	0.85
1:B:75:MET:HE1	1:B:105:VAL:HG11	1.58	0.83
1:A:237:ARG:HD2	1:A:242:GLU:OE1	1.77	0.82
1:B:176:TYR:CE1	1:D:176:TYR:CE1	2.68	0.82
1:D:193:LYS:HE3	1:D:194:ASN:H	1.49	0.78
1:B:22:ILE:HD12	1:B:23:SER:H	1.49	0.78
1:C:10:LYS:HB2	1:C:238:SER:OG	1.83	0.77
1:A:132:GLU:CD	1:A:133:LEU:H	1.87	0.75
1:A:126:ASN:OD1	1:A:133:LEU:HD21	1.89	0.72
1:A:204:ILE:HD11	1:D:185:ARG:CZ	2.20	0.72
1:A:132:GLU:CD	1:A:133:LEU:N	2.43	0.71
1:D:218:LEU:O	1:D:223:ARG:HD3	1.89	0.71
1:A:75:MET:HE1	1:A:105:VAL:CG1	2.22	0.70
1:C:218:LEU:O	1:C:223:ARG:HD3	1.92	0.70
1:C:165:VAL:O	1:C:231:LYS:HE3	1.91	0.69
1:B:75:MET:HE1	1:B:105:VAL:CG1	2.24	0.67
1:D:75:MET:HE1	1:D:105:VAL:CG1	2.23	0.66
1:C:6:THR:O	1:C:231:LYS:HE2	1.96	0.66
1:B:176:TYR:CE1	1:D:176:TYR:CD1	2.85	0.65
1:B:185:ARG:CZ	1:C:204:ILE:HD11	2.27	0.64
1:A:133:LEU:HG	1:A:135:LEU:HB2	1.80	0.63
1:B:58:ILE:HD11	1:B:133:LEU:HD22	1.82	0.62
1:C:324:ASN:ND2	1:D:298:ASP:H	1.99	0.60
1:B:139:HIS:HD2	1:B:152:GLU:OE2	1.85	0.59
1:B:176:TYR:CD1	1:D:176:TYR:CE1	2.90	0.59
1:C:139:HIS:HD2	1:C:152:GLU:OE2	1.86	0.58
1:B:139:HIS:HE1	1:B:169:THR:O	1.87	0.58
1:A:218:LEU:C	1:A:223:ARG:HH11	2.00	0.58
1:A:139:HIS:HD2	1:A:152:GLU:OE2	1.86	0.57
1:D:193:LYS:CE	1:D:194:ASN:H	2.16	0.57
1:D:139:HIS:HD2	1:D:152:GLU:OE2	1.88	0.57
1:B:56:ALA:O	1:B:57:GLY:C	2.41	0.56
1:B:204:ILE:HD11	1:C:185:ARG:CZ	2.36	0.56
1:A:2:CYS:N	1:A:170:ASN:HD21	2.04	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:ILE:HD11	1:D:185:ARG:NE	2.22	0.54
1:C:320:ILE:HD12	1:D:241:TYR:CE1	2.42	0.54
1:C:139:HIS:HE1	1:C:169:THR:O	1.91	0.54
1:C:2:CYS:O	1:C:223:ARG:NH2	2.41	0.53
1:D:139:HIS:HE1	1:D:169:THR:O	1.91	0.53
1:A:133:LEU:HD11	1:A:135:LEU:HD22	1.91	0.53
1:B:185:ARG:NE	1:C:204:ILE:HD11	2.24	0.52
1:A:139:HIS:HE1	1:A:169:THR:O	1.93	0.52
1:A:320:ILE:HD12	1:B:241:TYR:CE1	2.46	0.51
1:D:10:LYS:HB2	1:D:238:SER:HB2	1.92	0.50
1:C:241:TYR:CE1	1:D:320:ILE:HD12	2.47	0.50
1:A:241:TYR:CE1	1:B:320:ILE:HD12	2.47	0.49
1:A:99:PHE:CZ	1:A:135:LEU:HA	2.47	0.49
1:B:10:LYS:HB2	1:B:238:SER:HB2	1.94	0.49
1:D:264:GLY:O	1:D:266:GLU:N	2.44	0.48
1:A:132:GLU:OE1	1:A:133:LEU:N	2.46	0.48
1:D:2:CYS:O	1:D:223:ARG:NH2	2.46	0.48
1:B:139:HIS:CE1	1:B:169:THR:O	2.65	0.48
1:B:22:ILE:HD12	1:B:23:SER:N	2.25	0.48
1:A:185:ARG:NE	1:D:204:ILE:HD11	2.29	0.47
1:A:75:MET:CE	1:A:105:VAL:HB	2.45	0.47
1:A:139:HIS:CE1	1:A:169:THR:O	2.68	0.47
1:D:75:MET:CE	1:D:105:VAL:HB	2.44	0.47
1:C:139:HIS:CE1	1:C:169:THR:O	2.68	0.46
1:B:204:ILE:HD11	1:C:185:ARG:NE	2.31	0.46
1:A:229:PHE:CE2	1:B:260:LEU:HD12	2.51	0.45
1:B:75:MET:CE	1:B:105:VAL:HB	2.46	0.45
1:A:146:GLU:HG3	1:A:147:GLN:HB2	1.98	0.44
1:B:176:TYR:CE1	1:D:176:TYR:HE1	2.33	0.44
1:D:75:MET:HE2	1:D:75:MET:HB2	1.66	0.44
1:C:42:VAL:HG21	1:C:104:TRP:CZ3	2.52	0.44
1:B:277:CYS:SG	1:B:284:TYR:CD2	3.11	0.44
1:A:42:VAL:HG21	1:A:104:TRP:CZ3	2.53	0.44
1:B:266:GLU:O	1:B:266:GLU:HG2	2.18	0.43
1:D:193:LYS:HA	1:D:193:LYS:HD2	1.93	0.43
1:B:31:THR:HG21	1:B:37:PHE:CZ	2.54	0.43
1:D:75:MET:CE	1:D:105:VAL:CG1	2.96	0.42
1:D:139:HIS:CE1	1:D:169:THR:O	2.69	0.42
1:A:10:LYS:HB2	1:A:238:SER:HB2	2.01	0.42
1:B:75:MET:HB2	1:B:75:MET:HE2	1.68	0.42
1:A:185:ARG:CZ	1:D:204:ILE:HD11	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:TYR:HE1	1:D:176:TYR:CE1	2.32	0.42
1:D:184:TYR:HB3	1:D:187:LEU:HD22	2.00	0.42
1:A:277:CYS:SG	1:A:284:TYR:CD2	3.13	0.42
1:C:10:LYS:HA	1:C:10:LYS:HD2	1.83	0.41
1:A:126:ASN:OD1	1:A:133:LEU:CD2	2.62	0.41
1:C:31:THR:HG21	1:C:37:PHE:CZ	2.55	0.41
1:D:288:THR:HG21	1:D:316:GLU:HA	2.01	0.41
1:A:31:THR:HG21	1:A:37:PHE:CZ	2.55	0.41
1:B:24:TYR:N	1:B:26:GLU:OE2	2.48	0.41
1:C:36:LYS:HE3	1:C:46:ASP:OD1	2.21	0.41
1:B:36:LYS:HE3	1:B:46:ASP:OD1	2.21	0.41
1:A:75:MET:CE	1:A:105:VAL:CG1	2.96	0.40
1:A:40:ARG:HD2	1:A:129:PHE:HB2	2.04	0.40

All (2) 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)
2:A:405:HOH:O	2:C:401:HOH:O[1_655]	1.24	0.96
1:A:24:TYR:OH	1:C:20:TYR:OH[1_655]	2.15	0.05

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	321/331 (97%)	308 (96%)	13 (4%)	0	100	100
1	B	321/331 (97%)	309 (96%)	11 (3%)	1 (0%)	41	37
1	C	321/331 (97%)	314 (98%)	7 (2%)	0	100	100
1	D	321/331 (97%)	306 (95%)	15 (5%)	0	100	100
All	All	1284/1324 (97%)	1237 (96%)	46 (4%)	1 (0%)	51	49

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	265	ASP

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	290/298 (97%)	283 (98%)	7 (2%)	49	51
1	B	290/298 (97%)	287 (99%)	3 (1%)	76	81
1	C	290/298 (97%)	282 (97%)	8 (3%)	43	44
1	D	290/298 (97%)	281 (97%)	9 (3%)	40	40
All	All	1160/1192 (97%)	1133 (98%)	27 (2%)	50	53

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	16	ARG
1	A	119	LEU
1	A	130	SER
1	A	132	GLU
1	A	135	LEU
1	A	139	HIS
1	B	16	ARG
1	B	139	HIS
1	B	193	LYS
1	C	16	ARG
1	C	24	TYR
1	C	119	LEU
1	C	139	HIS
1	C	193	LYS
1	C	201	GLU
1	C	231	LYS
1	C	237	ARG
1	D	16	ARG

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Mol	Chain	Res	Type
1	D	20	TYR
1	D	58	ILE
1	D	139	HIS
1	D	156	GLU
1	D	187	LEU
1	D	193	LYS
1	D	201	GLU
1	D	288	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	HIS
1	A	183	ASN
1	B	25	ASN
1	B	139	HIS
1	C	139	HIS
1	C	324	ASN
1	D	139	HIS
1	D	183	ASN

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 ⓘ

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, 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	323/331 (97%)	-0.09	10 (3%) 49 48	9, 16, 38, 120	0
1	B	323/331 (97%)	-0.25	7 (2%) 62 60	10, 15, 31, 68	0
1	C	323/331 (97%)	-0.21	6 (1%) 66 65	10, 15, 32, 63	0
1	D	323/331 (97%)	-0.15	7 (2%) 62 60	10, 16, 34, 61	0
All	All	1292/1324 (97%)	-0.17	30 (2%) 60 59	9, 16, 34, 120	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	129	PHE	13.2
1	A	130	SER	11.6
1	A	131	ASP	7.3
1	A	133	LEU	6.9
1	D	59	ALA	6.4
1	A	128	ASN	6.2
1	B	265	ASP	5.2
1	A	132	GLU	4.8
1	C	133	LEU	4.4
1	D	266	GLU	4.1
1	A	134	PRO	4.1
1	D	265	ASP	3.9
1	C	24	TYR	3.8
1	C	60	ASP	3.8
1	B	264	GLY	3.7
1	A	60	ASP	3.5
1	B	24	TYR	3.2
1	A	24	TYR	3.2
1	D	264	GLY	2.8
1	B	266	GLU	2.7
1	A	58	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	131	ASP	2.5
1	C	132	GLU	2.3
1	C	58	ILE	2.3
1	D	58	ILE	2.3
1	B	198	ASP	2.2
1	B	47	HIS	2.2
1	B	104	TRP	2.2
1	D	60	ASP	2.1
1	C	131	ASP	2.1

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.