



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:00 AM UTC

PDB ID : 4WL3 / pdb_00004wl3
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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

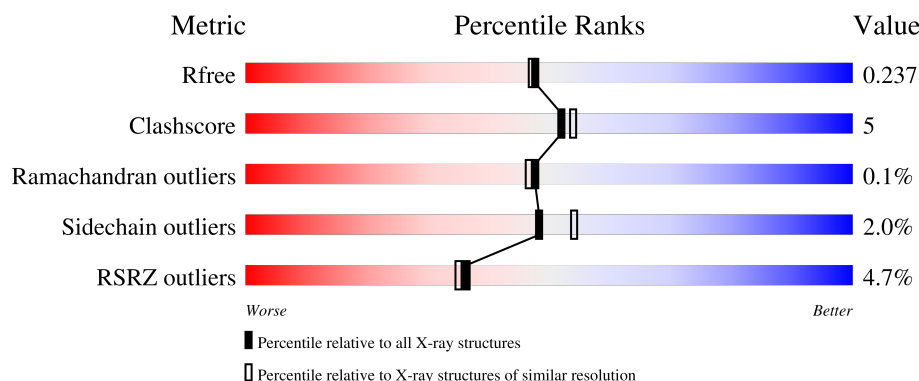
1 Overall quality at a glance

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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 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 $\leq 5\%$. 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	6% 84% 12% ..
1	B	331	3% 85% 10% ..
1	C	331	5% 88% 8% ..
1	D	331	4% 88% 7% ...

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

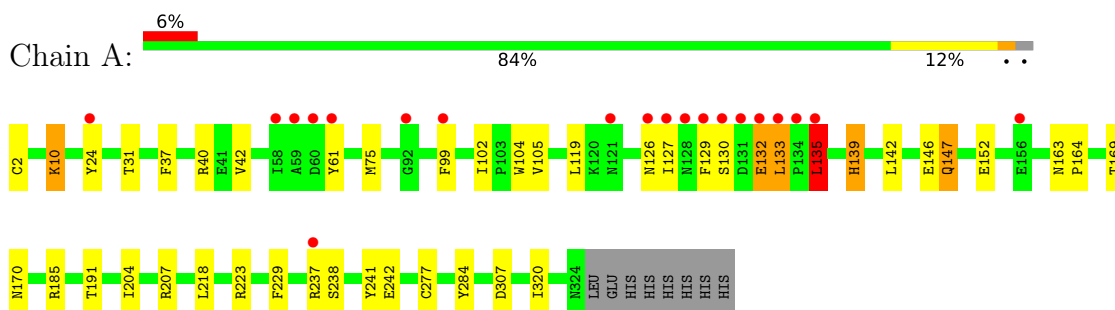
- Molecule 2 is water.

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

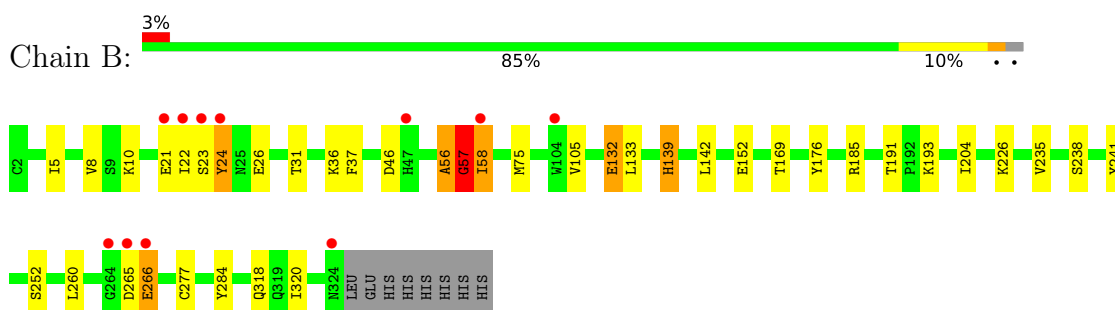
3 Residue-property plots [i](#)

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.

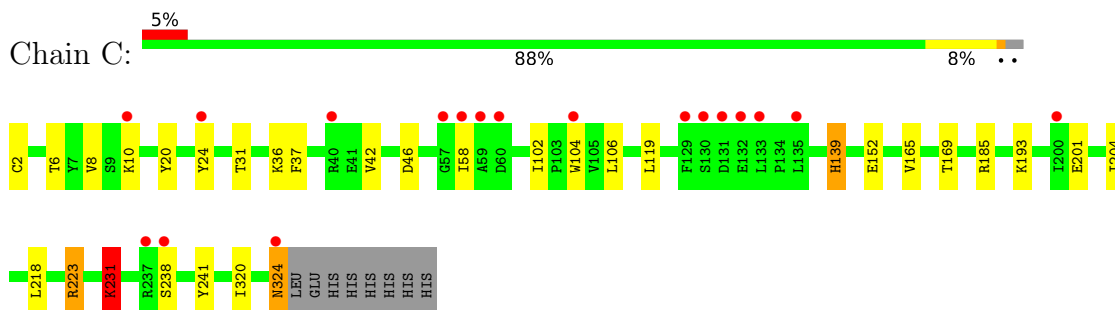
- Molecule 1: Bile salt hydrolase



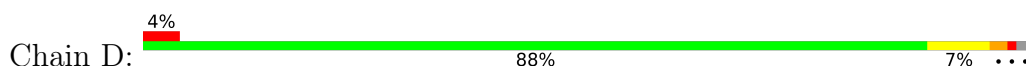
- Molecule 1: Bile salt hydrolase

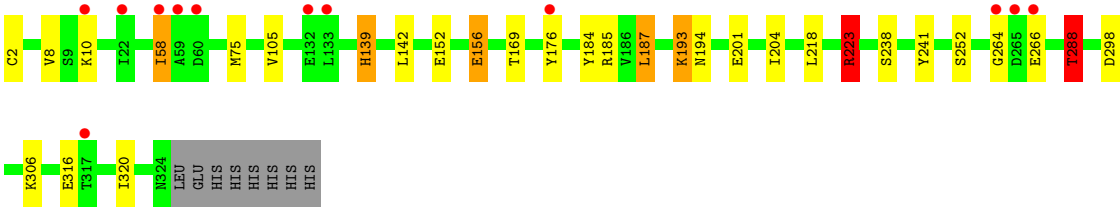


- Molecule 1: Bile salt hydrolase



- Molecule 1: Bile salt hydrolase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.20Å 131.62Å 86.72Å 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) 95.3 (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 2.01Å)	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 (4.74%)	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)	0.40 , 30.1	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	1.24	5/2656 (0.2%)	1.06	7/3601 (0.2%)
1	B	1.18	4/2656 (0.2%)	1.10	12/3601 (0.3%)
1	C	1.18	4/2656 (0.2%)	1.06	5/3601 (0.1%)
1	D	1.18	2/2656 (0.1%)	1.07	6/3601 (0.2%)
All	All	1.20	15/10624 (0.1%)	1.07	30/14404 (0.2%)

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 (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	57	GLY	N-CA	7.83	1.57	1.45
1	A	102	ILE	CA-CB	7.60	1.58	1.54
1	D	223	ARG	CD-NE	-7.22	1.36	1.46
1	C	223	ARG	CD-NE	-6.55	1.37	1.46
1	A	135	LEU	N-CA	5.98	1.53	1.46
1	A	164	PRO	C-O	-5.78	1.16	1.24
1	C	102	ILE	CA-CB	5.31	1.57	1.54
1	D	252	SER	C-O	-5.28	1.17	1.24
1	B	252	SER	C-O	-5.26	1.17	1.24
1	C	231	LYS	N-CA	5.19	1.52	1.46
1	B	191	THR	C-O	-5.13	1.17	1.24
1	A	163	ASN	C-O	-5.11	1.18	1.24
1	B	5	ILE	N-CA	-5.10	1.40	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	106	LEU	C-O	-5.09	1.17	1.24
1	A	191	THR	C-O	-5.08	1.17	1.24

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	223	ARG	NE-CZ-NH2	-11.52	108.83	119.20
1	D	223	ARG	NE-CZ-NH2	-11.20	109.12	119.20
1	C	223	ARG	NE-CZ-NH1	9.60	131.10	121.50
1	D	288	THR	N-CA-CB	-9.04	97.06	111.43
1	D	223	ARG	CD-NE-CZ	8.73	136.62	124.40
1	B	56	ALA	CA-C-N	-8.41	103.89	121.06
1	B	56	ALA	C-N-CA	-8.41	103.89	121.06
1	A	133	LEU	CA-C-N	-8.38	111.10	119.56
1	A	133	LEU	C-N-CA	-8.38	111.10	119.56
1	D	223	ARG	NE-CZ-NH1	8.22	129.72	121.50
1	A	147	GLN	N-CA-CB	-7.86	99.39	111.56
1	B	226	LYS	CG-CD-CE	7.54	128.63	111.30
1	B	57	GLY	CA-C-O	7.23	128.06	119.82
1	A	135	LEU	N-CA-CB	6.93	119.86	109.19
1	B	266	GLU	N-CA-C	-6.68	105.02	112.57
1	C	324	ASN	N-CA-CB	-6.61	99.27	110.50
1	B	58	ILE	N-CA-C	-6.57	106.07	111.91
1	C	223	ARG	CD-NE-CZ	6.54	133.56	124.40
1	A	61	TYR	N-CA-CB	-6.33	102.37	110.03
1	B	56	ALA	O-C-N	-6.31	115.22	122.03
1	B	24	TYR	CB-CA-C	-5.66	103.06	111.72
1	B	8	VAL	CB-CA-C	-5.64	102.22	110.62
1	A	10	LYS	N-CA-C	-5.56	105.54	112.93
1	B	318	GLN	CB-CA-C	-5.49	101.33	109.90
1	D	8	VAL	CB-CA-C	-5.29	102.73	110.62
1	D	306	LYS	CD-CE-NZ	5.29	128.82	111.90
1	C	8	VAL	CB-CA-C	-5.25	102.82	110.81
1	B	132	GLU	CB-CA-C	-5.14	101.12	110.63
1	A	307	ASP	CB-CA-C	-5.12	99.97	109.29
1	B	58	ILE	N-CA-CB	-5.03	106.50	112.39

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	127	ILE	Peptide

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Mol	Chain	Res	Type	Group
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	37	1
1	B	2597	0	2504	34	0
1	C	2597	0	2504	19	1
1	D	2597	0	2504	33	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	103	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 5.

All (103) 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:D:75:MET:HE1	1:D:105:VAL:HB	1.59	0.84
1:A:75:MET:HE1	1:A:105:VAL:CG1	2.07	0.83
1:D:75:MET:HE1	1:D:105:VAL:CG1	2.08	0.83
1:A:237:ARG:HD2	1:A:242:GLU:OE1	1.77	0.82
1:D:75:MET:HE1	1:D:105:VAL:CB	2.09	0.82
1:B:176:TYR:CE1	1:D:176:TYR:CE1	2.68	0.82
1:A:132:GLU:CD	1:A:133:LEU:H	1.87	0.81
1:B:75:MET:HE1	1:B:105:VAL:CB	2.11	0.81
1:B:75:MET:HE1	1:B:105:VAL:CG1	2.10	0.80
1:A:75:MET:HE1	1:A:105:VAL:HB	1.62	0.80
1:A:75:MET:HE1	1:A:105:VAL:CB	2.11	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:MET:HE1	1:B:105:VAL:HB	1.61	0.80
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:N	2.43	0.76
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:D:218:LEU:O	1:D:223:ARG:HD3	1.89	0.71
1:C:218:LEU:O	1:C:223:ARG:HD3	1.92	0.70
1:A:75:MET:HE1	1:A:105:VAL:HG11	1.74	0.69
1:C:165:VAL:O	1:C:231:LYS:HE3	1.91	0.69
1:D:75:MET:HE1	1:D:105:VAL:HG11	1.75	0.68
1:B:75:MET:HE1	1:B:105:VAL:HG11	1.78	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:A:218:LEU:C	1:A:223:ARG:HH11	2.00	0.62
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:56:ALA:O	1:B:57:GLY:C	2.41	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: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: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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:PHE:CZ	1:A:135:LEU:HA	2.47	0.49
1:A:241:TYR:CE1	1:B:320:ILE:HD12	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:HE2	1:A:142:LEU:CD2	2.44	0.47
1:A:139:HIS:CE1	1:A:169:THR:O	2.68	0.47
1:A:75:MET:HE2	1:A:142:LEU:HG	1.98	0.46
1:C:139:HIS:CE1	1:C:169:THR:O	2.68	0.46
1:D:75:MET:CE	1:D:105:VAL:HG11	2.44	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:D:75:MET:HE2	1:D:142:LEU:HG	1.99	0.45
1:B:75:MET:HE2	1:B:142:LEU:CD2	2.47	0.45
1:D:75:MET:HE2	1:D:142:LEU:CD2	2.47	0.45
1:A:75:MET:HE3	1:A:75:MET:HB2	1.73	0.44
1:A:146:GLU:HG3	1:A:147:GLN:HB2	1.98	0.44
1:A:75:MET:CE	1:A:105:VAL:HG11	2.45	0.44
1:B:176:TYR:CE1	1:D:176:TYR:HE1	2.33	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:HD2	1:D:193:LYS:HA	1.93	0.43
1:B:75:MET:HE2	1:B:142:LEU:HG	1.99	0.43
1:D:156:GLU:H	1:D:156:GLU:CD	2.26	0.43
1:B:75:MET:CE	1:B:105:VAL:HG11	2.47	0.43
1:B:31:THR:HG21	1:B:37:PHE:CZ	2.54	0.43
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:A:185:ARG:CZ	1:D:204:ILE:HD11	2.48	0.42
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:58:ILE:HD12	1:D:58:ILE:C	2.46	0.41
1:B:75:MET:HE3	1:B:75:MET:HB2	1.73	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: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%)	36	35
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%)	48	46

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%)	285 (98%)	5 (2%)	53	60
1	B	290/298 (97%)	286 (99%)	4 (1%)	59	66
1	C	290/298 (97%)	283 (98%)	7 (2%)	43	47
1	D	290/298 (97%)	283 (98%)	7 (2%)	43	47
All	All	1160/1192 (97%)	1137 (98%)	23 (2%)	48	54

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	LEU
1	A	130	SER
1	A	132	GLU
1	A	135	LEU
1	A	139	HIS
1	B	132	GLU
1	B	139	HIS
1	B	193	LYS
1	B	235	VAL
1	C	24	TYR
1	C	58	ILE
1	C	119	LEU
1	C	139	HIS
1	C	193	LYS
1	C	201	GLU
1	C	231	LYS
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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	HIS
1	A	183	ASN
1	A	300	ASN
1	B	25	ASN
1	B	139	HIS
1	C	139	HIS
1	C	183	ASN
1	C	293	GLN
1	C	324	ASN
1	D	139	HIS
1	D	147	GLN
1	D	183	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 oligosaccharides 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, 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%)	0.16	20 (6%)	26 25	9, 16, 38, 120	0
1	B	323/331 (97%)	-0.06	11 (3%)	48 47	10, 15, 31, 68	0
1	C	323/331 (97%)	0.03	18 (5%)	30 29	10, 15, 32, 63	0
1	D	323/331 (97%)	0.01	12 (3%)	45 44	10, 16, 34, 61	0
All	All	1292/1324 (97%)	0.04	61 (4%)	36 35	9, 16, 34, 120	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	129	PHE	10.4
1	A	133	LEU	8.5
1	D	59	ALA	8.3
1	A	128	ASN	7.8
1	A	134	PRO	7.6
1	C	24	TYR	7.2
1	A	131	ASP	6.5
1	A	130	SER	6.3
1	A	132	GLU	6.1
1	D	58	ILE	5.7
1	B	24	TYR	5.4
1	C	133	LEU	5.2
1	A	24	TYR	5.0
1	C	132	GLU	4.9
1	A	127	ILE	4.2
1	B	104	TRP	4.0
1	D	60	ASP	3.8
1	C	58	ILE	3.5
1	B	266	GLU	3.4
1	B	265	ASP	3.3
1	A	58	ILE	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	135	LEU	3.3
1	B	264	GLY	3.2
1	D	132	GLU	3.2
1	C	60	ASP	3.1
1	A	60	ASP	3.1
1	A	92	GLY	3.1
1	D	176	TYR	3.0
1	C	237	ARG	3.0
1	D	264	GLY	2.8
1	A	156	GLU	2.8
1	D	266	GLU	2.8
1	C	104	TRP	2.8
1	B	324	ASN	2.7
1	B	21	GLU	2.7
1	C	40	ARG	2.7
1	A	121	ASN	2.6
1	C	129	PHE	2.6
1	D	265	ASP	2.6
1	A	61	TYR	2.6
1	C	324	ASN	2.6
1	A	99	PHE	2.5
1	D	22	ILE	2.5
1	C	59	ALA	2.5
1	D	133	LEU	2.4
1	C	131	ASP	2.4
1	B	23	SER	2.4
1	A	126	ASN	2.4
1	D	317	THR	2.3
1	B	58	ILE	2.3
1	C	200	ILE	2.2
1	C	57	GLY	2.2
1	C	135	LEU	2.2
1	C	130	SER	2.2
1	B	47	HIS	2.1
1	A	237	ARG	2.1
1	C	10	LYS	2.1
1	D	10	LYS	2.1
1	A	59	ALA	2.1
1	B	22	ILE	2.1
1	C	238	SER	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 oligosaccharides 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.