



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:33 PM UTC

PDB ID : 6LD6 / pdb_00006ld6
Title : Structure of Bifidobacterium dentium beta-glucuronidase
Authors : Lin, H.-Y.; Hsieh, T.-J.; Lin, C.-H.
Deposited on : 2019-11-20
Resolution : 2.20 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

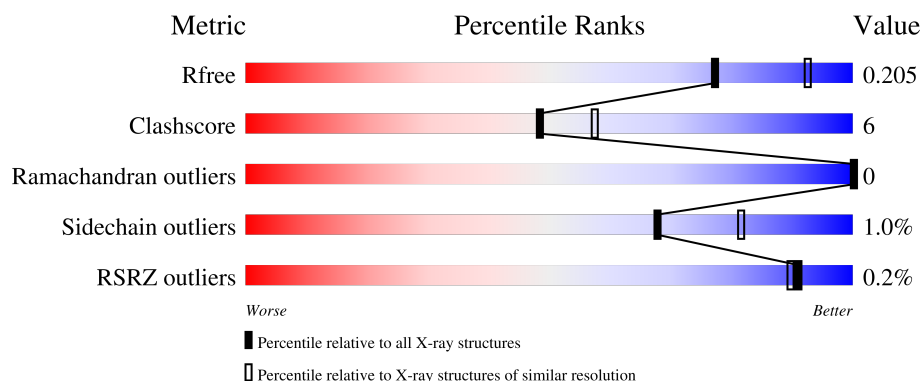
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	670	 82% 9% • 7%
1	B	670	 83% 9% • 7%
1	C	670	 83% 9% • 7%
1	D	670	 78% 14% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 21571 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 LacZ1 Beta-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	620	Total	C	N	O	S	0	0	0
			4909	3113	856	921	19			
1	B	620	Total	C	N	O	S	0	0	0
			4909	3113	856	921	19			
1	C	620	Total	C	N	O	S	0	0	0
			4909	3113	856	921	19			
1	D	620	Total	C	N	O	S	0	0	0
			4909	3113	856	921	19			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ASN	-	expression tag	UNP D2Q7B1
A	0	GLY	-	expression tag	UNP D2Q7B1
B	-1	ASN	-	expression tag	UNP D2Q7B1
B	0	GLY	-	expression tag	UNP D2Q7B1
C	-1	ASN	-	expression tag	UNP D2Q7B1
C	0	GLY	-	expression tag	UNP D2Q7B1
D	-1	ASN	-	expression tag	UNP D2Q7B1
D	0	GLY	-	expression tag	UNP D2Q7B1

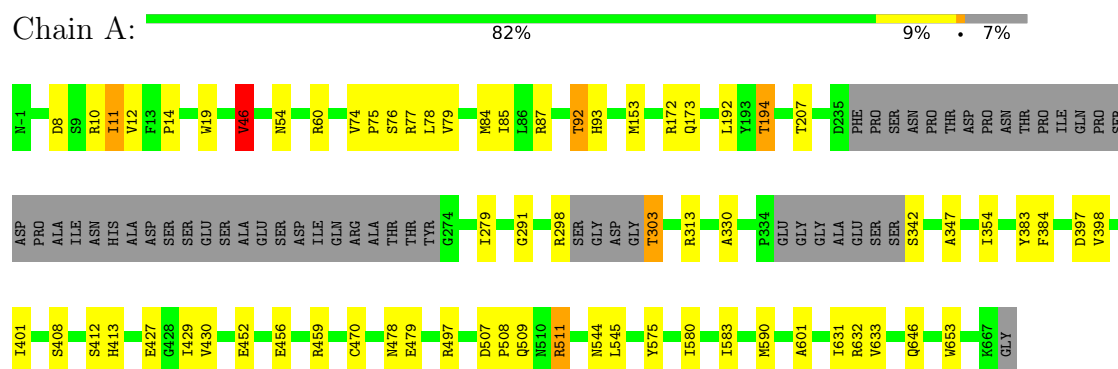
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	527	Total	O	0	0
			527	527		
2	B	455	Total	O	0	0
			455	455		
2	C	521	Total	O	0	0
			521	521		
2	D	432	Total	O	0	0
			432	432		

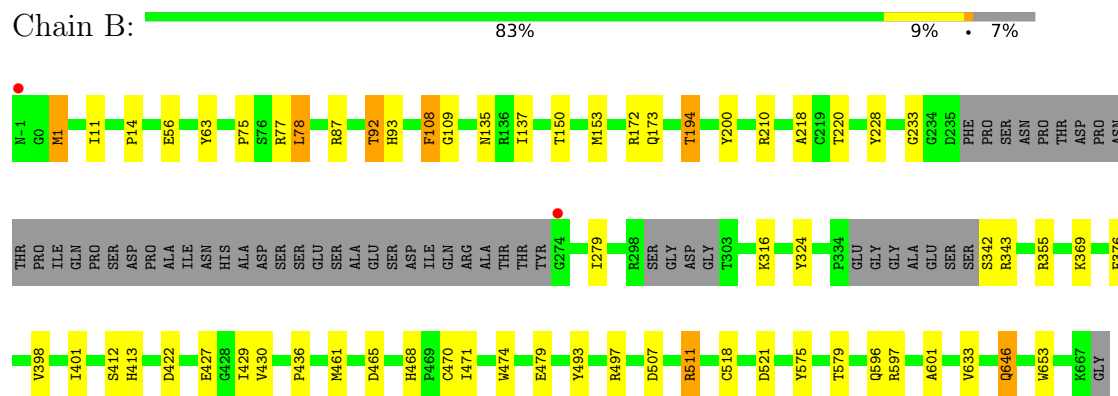
3 Residue-property plots

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.

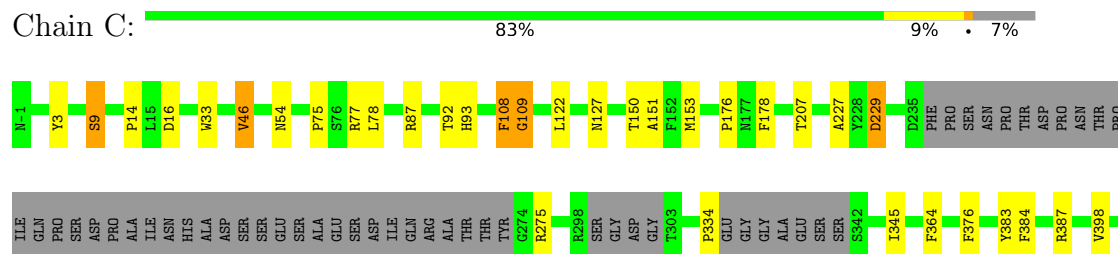
• Molecule 1: LacZ1 Beta-galactosidase



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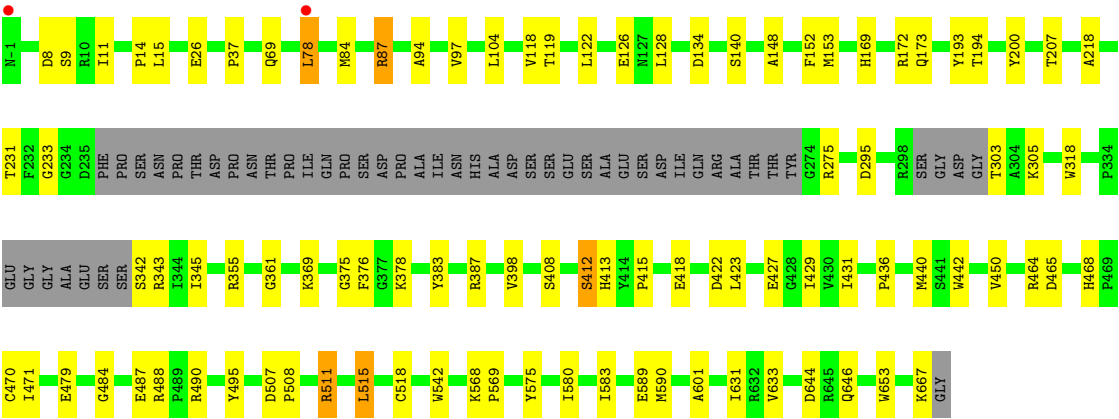
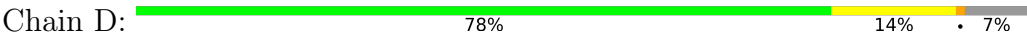


• Molecule 1: LacZ1 Beta-galactosidase





● Molecule 1: LacZ1 Beta-galactosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	92.58Å 104.90Å 160.62Å 90.00° 91.16° 90.00°	Depositor
Resolution (Å)	29.60 – 2.20 29.60 – 2.20	Depositor EDS
% Data completeness (in resolution range)	94.9 (29.60-2.20) 94.7 (29.60-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.41 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.10_2155	Depositor
R, R_{free}	0.151 , 0.205 0.152 , 0.205	Depositor DCC
R_{free} test set	2012 reflections (1.30%)	wwPDB-VP
Wilson B-factor (Å ²)	22.3	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.045 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	21571	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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	1.12	5/5040 (0.1%)	1.07	9/6857 (0.1%)
1	B	1.04	4/5040 (0.1%)	1.04	7/6857 (0.1%)
1	C	1.12	9/5040 (0.2%)	1.12	13/6857 (0.2%)
1	D	1.00	1/5040 (0.0%)	1.05	12/6857 (0.2%)
All	All	1.07	19/20160 (0.1%)	1.07	41/27428 (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

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	109	GLY	N-CA	10.03	1.56	1.45
1	A	511	ARG	CD-NE	-8.44	1.34	1.46
1	B	78	LEU	CG-CD2	7.97	1.78	1.52
1	A	46	VAL	CA-CB	7.44	1.63	1.54
1	C	109	GLY	CA-C	7.37	1.59	1.52
1	C	511	ARG	CD-NE	-6.71	1.36	1.46
1	C	108	PHE	C-N	-6.58	1.27	1.33
1	C	46	VAL	CA-CB	6.41	1.62	1.54
1	B	92	THR	CA-C	6.33	1.60	1.52
1	A	384	PHE	N-CA	6.26	1.54	1.45
1	B	1	MET	CA-CB	5.80	1.63	1.53
1	A	60	ARG	CZ-NH1	5.46	1.40	1.32
1	C	632	ARG	CZ-NH1	5.27	1.40	1.32
1	C	229	ASP	CB-CG	-5.22	1.39	1.52
1	B	511	ARG	CD-NE	-5.14	1.39	1.46
1	D	78	LEU	CB-CG	-5.10	1.43	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	92	THR	CB-CG2	5.08	1.69	1.52
1	C	412	SER	CA-C	5.05	1.59	1.52
1	C	384	PHE	N-CA	5.03	1.52	1.45

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	108	PHE	CA-C-N	-18.84	101.83	119.92
1	C	108	PHE	C-N-CA	-18.84	101.83	119.92
1	C	109	GLY	N-CA-C	12.22	125.66	111.36
1	B	78	LEU	CB-CG-CD1	-11.48	76.26	110.70
1	D	78	LEU	CB-CG-CD1	-11.35	76.65	110.70
1	C	383	TYR	CA-C-N	-11.34	104.43	122.24
1	C	383	TYR	C-N-CA	-11.34	104.43	122.24
1	A	383	TYR	CA-C-N	-9.78	106.89	122.24
1	A	383	TYR	C-N-CA	-9.78	106.89	122.24
1	D	383	TYR	CA-C-N	-9.25	106.93	122.11
1	D	383	TYR	C-N-CA	-9.25	106.93	122.11
1	A	511	ARG	NE-CZ-NH2	-9.18	110.94	119.20
1	A	87	ARG	CG-CD-NE	-8.87	92.49	112.00
1	C	87	ARG	CG-CD-NE	-7.87	94.69	112.00
1	C	87	ARG	NE-CZ-NH2	-7.80	112.18	119.20
1	C	511	ARG	NE-CZ-NH2	-7.73	112.24	119.20
1	C	108	PHE	O-C-N	-7.22	112.21	122.03
1	B	87	ARG	NE-CZ-NH2	-7.06	112.85	119.20
1	B	511	ARG	CB-CG-CD	-6.64	96.02	111.30
1	D	87	ARG	NE-CZ-NH2	-6.59	113.27	119.20
1	A	153	MET	CB-CG-SD	-6.55	93.05	112.70
1	D	511	ARG	CB-CG-CD	-6.12	97.24	111.30
1	B	87	ARG	CG-CD-NE	-5.89	99.05	112.00
1	A	87	ARG	NE-CZ-NH2	-5.83	113.95	119.20
1	B	108	PHE	O-C-N	-5.71	114.26	122.03
1	C	153	MET	CB-CG-SD	-5.63	95.80	112.70
1	C	33	TRP	CA-C-N	-5.57	112.31	122.38
1	C	33	TRP	C-N-CA	-5.57	112.31	122.38
1	B	78	LEU	CA-CB-CG	-5.54	96.91	116.30
1	B	646	GLN	CA-CB-CG	-5.53	103.05	114.10
1	D	148	ALA	CA-C-N	-5.39	117.28	121.86
1	D	148	ALA	C-N-CA	-5.39	117.28	121.86
1	A	511	ARG	NE-CZ-NH1	5.35	126.85	121.50
1	D	484	GLY	CA-C-N	-5.26	112.89	123.09
1	D	484	GLY	C-N-CA	-5.26	112.89	123.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	383	TYR	N-CA-C	5.23	117.39	111.11
1	A	590	MET	CA-C-N	5.14	131.35	121.54
1	A	590	MET	C-N-CA	5.14	131.35	121.54
1	D	193	TYR	CA-C-N	5.04	130.23	122.93
1	D	193	TYR	C-N-CA	5.04	130.23	122.93
1	C	3	TYR	CA-CB-CG	-5.01	104.89	113.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	303	THR	Peptide

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	4909	0	4659	50	0
1	B	4909	0	4659	67	0
1	C	4909	0	4659	41	0
1	D	4909	0	4659	70	0
2	A	527	0	0	11	1
2	B	455	0	0	16	0
2	C	521	0	0	8	1
2	D	432	0	0	16	0
All	All	21571	0	18636	211	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 6.

All (211) 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:B:78:LEU:CD2	1:B:78:LEU:CG	1.78	1.58
1:B:78:LEU:HA	1:B:78:LEU:HD23	1.33	1.10

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:ARG:NH2	2:B:701:HOH:O	1.86	1.08
1:D:152:PHE:CD2	1:D:153:MET:HE2	1.97	0.98
1:B:78:LEU:CD2	1:B:78:LEU:HA	1.97	0.95
1:B:78:LEU:CD2	1:B:78:LEU:CA	2.47	0.93
1:B:78:LEU:CD2	1:B:78:LEU:CB	2.48	0.92
1:A:77:ARG:HD2	2:C:1109:HOH:O	1.76	0.86
1:A:646:GLN:OE1	2:A:701:HOH:O	1.94	0.85
1:A:459:ARG:NH1	2:A:702:HOH:O	2.03	0.82
1:B:78:LEU:HD23	1:B:78:LEU:CA	2.07	0.82
1:D:231:THR:OG1	1:D:305:LYS:NZ	2.14	0.79
1:A:497:ARG:NH1	2:A:703:HOH:O	2.07	0.76
1:D:172:ARG:NH2	2:D:701:HOH:O	2.10	0.75
1:C:497:ARG:NH1	2:C:701:HOH:O	2.09	0.74
1:B:172:ARG:NH1	2:B:708:HOH:O	2.21	0.73
1:B:342:SER:N	2:B:706:HOH:O	2.21	0.73
1:D:342:SER:N	2:D:707:HOH:O	2.22	0.72
1:B:14:PRO:HG2	1:D:14:PRO:HG2	1.71	0.71
1:D:218:ALA:HB2	1:D:369:LYS:HE2	1.72	0.71
1:B:355:ARG:NH2	1:B:511:ARG:HH21	1.88	0.71
1:B:343:ARG:NH2	2:B:706:HOH:O	2.24	0.70
1:D:152:PHE:CD2	1:D:153:MET:CE	2.75	0.70
1:B:56:GLU:OE1	2:B:702:HOH:O	2.09	0.69
1:D:152:PHE:CG	1:D:153:MET:HE2	2.27	0.68
1:A:298:ARG:NH2	1:A:303:THR:HB	2.09	0.67
1:B:646:GLN:OE1	2:B:703:HOH:O	2.14	0.65
1:A:580:ILE:HB	1:A:583:ILE:HD12	1.78	0.65
1:D:303:THR:N	2:D:714:HOH:O	2.30	0.65
1:B:633:VAL:HG21	1:C:633:VAL:HG21	1.80	0.63
1:D:87:ARG:NH2	1:D:418:GLU:OE2	2.32	0.63
1:B:210:ARG:NH2	2:B:709:HOH:O	2.31	0.63
1:A:75:PRO:HD2	1:A:78:LEU:HD22	1.80	0.62
1:C:376:PHE:HZ	1:C:401:ILE:HG12	1.64	0.62
1:A:298:ARG:HH12	1:A:303:THR:N	1.98	0.61
1:D:490:ARG:NH1	2:D:713:HOH:O	2.29	0.61
1:B:398:VAL:HG11	1:B:427:GLU:HG3	1.83	0.61
1:B:93:HIS:O	1:B:108:PHE:O	2.19	0.61
1:A:452:GLU:OE2	2:A:704:HOH:O	2.16	0.61
1:D:169:HIS:ND1	2:D:716:HOH:O	2.31	0.61
1:B:343:ARG:NH1	2:B:706:HOH:O	2.33	0.60
1:A:298:ARG:CZ	1:A:303:THR:HB	2.32	0.59
1:B:479:GLU:HB3	1:B:518:CYS:HB3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:78:LEU:HD22	1:D:78:LEU:CD2	2.33	0.58
1:C:580:ILE:HB	1:C:583:ILE:HD12	1.85	0.58
1:A:173:GLN:HB3	2:A:818:HOH:O	2.04	0.58
1:B:597:ARG:NH2	2:B:713:HOH:O	2.37	0.57
1:C:398:VAL:HG11	1:C:427:GLU:HG3	1.86	0.57
1:B:465:ASP:HB3	1:B:471:ILE:CD1	2.35	0.56
1:B:78:LEU:CD2	1:D:78:LEU:HD22	2.35	0.56
1:B:220:THR:HG22	1:B:316:LYS:HE2	1.87	0.56
1:B:78:LEU:CD2	1:B:78:LEU:CD1	2.76	0.56
1:B:78:LEU:HD22	1:D:78:LEU:HD22	1.88	0.56
1:A:78:LEU:HG	1:C:78:LEU:HD21	1.88	0.55
1:A:303:THR:N	2:A:717:HOH:O	2.39	0.55
1:D:84:MET:HG3	1:D:122:LEU:HD12	1.87	0.55
1:B:343:ARG:CZ	2:B:706:HOH:O	2.55	0.55
1:B:355:ARG:HH22	1:B:511:ARG:NH2	2.05	0.55
1:D:507:ASP:OD2	1:D:511:ARG:HD2	2.07	0.55
1:B:150:THR:HB	2:B:911:HOH:O	2.07	0.55
1:D:295:ASP:OD2	2:D:702:HOH:O	2.18	0.54
1:B:461:MET:HE3	1:B:474:TRP:CE3	2.43	0.54
1:D:207:THR:HG21	1:D:508:PRO:HB2	1.90	0.54
1:B:376:PHE:HZ	1:B:401:ILE:HG12	1.73	0.53
1:A:77:ARG:HH22	1:C:9:SER:HA	1.72	0.53
1:A:14:PRO:HG2	1:C:14:PRO:HG2	1.90	0.53
1:A:194:THR:HG21	1:C:77:ARG:HH22	1.73	0.53
1:B:56:GLU:OE2	2:B:704:HOH:O	2.18	0.52
1:B:493:TYR:CZ	1:B:497:ARG:HG3	2.44	0.52
1:C:644:ASP:OD1	2:C:702:HOH:O	2.19	0.52
1:A:478:ASN:ND2	1:A:479:GLU:HG3	2.25	0.51
1:C:601:ALA:HA	1:C:653:TRP:CH2	2.46	0.51
1:D:303:THR:HG23	1:D:303:THR:O	2.11	0.51
1:A:78:LEU:HG	1:C:78:LEU:CD2	2.41	0.51
1:C:465:ASP:HB3	1:C:471:ILE:CD1	2.41	0.50
1:B:422:ASP:OD1	1:B:468:HIS:HE1	1.95	0.50
1:A:76:SER:HB2	2:A:1066:HOH:O	2.12	0.50
1:D:580:ILE:HB	1:D:583:ILE:HD12	1.94	0.50
1:B:601:ALA:HA	1:B:653:TRP:CH2	2.47	0.50
1:D:87:ARG:HH22	1:D:418:GLU:CD	2.19	0.50
1:D:378:LYS:NZ	2:D:718:HOH:O	2.39	0.49
1:B:108:PHE:O	1:B:109:GLY:C	2.53	0.49
1:B:355:ARG:NH2	1:B:511:ARG:NH2	2.58	0.49
1:A:398:VAL:HG11	1:A:427:GLU:HG3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:343:ARG:NH2	1:D:345:ILE:HD11	2.28	0.48
1:A:11:ILE:HD12	1:A:12:VAL:H	1.78	0.48
1:A:79:VAL:HG12	1:A:84:MET:HE2	1.96	0.48
1:D:200:TYR:CZ	1:D:233:GLY:HA3	2.48	0.48
1:A:54:ASN:C	2:A:706:HOH:O	2.57	0.48
1:A:544:ASN:O	1:A:545:LEU:C	2.55	0.48
1:C:497:ARG:NH2	2:C:709:HOH:O	2.34	0.48
1:D:487:GLU:OE1	2:D:704:HOH:O	2.20	0.48
1:A:509:GLN:HB3	1:A:511:ARG:HD2	1.95	0.48
1:C:176:PRO:HB3	1:C:178:PHE:CZ	2.49	0.48
1:A:10:ARG:HA	1:A:194:THR:HG22	1.96	0.48
1:A:298:ARG:HH22	1:A:303:THR:HB	1.76	0.48
1:D:415:PRO:HD3	1:D:436:PRO:HD3	1.96	0.48
1:B:172:ARG:CZ	2:B:701:HOH:O	2.48	0.47
1:C:364:PHE:HB2	1:C:614:ILE:HD13	1.95	0.47
1:A:298:ARG:NH1	1:A:303:THR:HB	2.30	0.47
1:C:429:ILE:O	1:C:470:CYS:HB2	2.15	0.47
1:D:126:GLU:HG2	2:D:715:HOH:O	2.14	0.47
1:B:412:SER:HA	1:B:413:HIS:HA	1.75	0.47
1:B:633:VAL:HG21	1:C:633:VAL:CG2	2.45	0.47
1:C:649:MET:HE3	1:C:649:MET:HB2	1.64	0.47
1:A:429:ILE:O	1:A:470:CYS:HB2	2.15	0.46
1:C:108:PHE:O	1:C:109:GLY:C	2.50	0.46
1:D:361:GLY:O	1:D:569:PRO:HD3	2.15	0.46
1:A:408:SER:HA	1:A:430:VAL:O	2.15	0.46
1:B:75:PRO:HA	1:D:8:ASP:HA	1.96	0.46
1:D:26:GLU:HG3	2:D:898:HOH:O	2.14	0.46
1:A:452:GLU:O	1:A:456:GLU:HG3	2.16	0.46
1:C:207:THR:OG1	1:C:227:ALA:HB3	2.15	0.46
1:A:354:ILE:HD13	1:A:354:ILE:HG21	1.65	0.46
1:A:633:VAL:HG21	1:D:633:VAL:HG21	1.98	0.46
1:B:77:ARG:HG3	1:B:78:LEU:HD21	1.98	0.46
1:B:479:GLU:CB	1:B:518:CYS:HB3	2.46	0.46
1:C:75:PRO:HB2	1:C:78:LEU:HD12	1.98	0.45
1:C:412:SER:HA	1:C:413:HIS:HA	1.76	0.45
1:C:92:THR:HA	1:C:93:HIS:HA	1.72	0.45
1:D:398:VAL:HG11	1:D:427:GLU:HG3	1.99	0.45
1:D:631:ILE:HD13	1:D:631:ILE:HG21	1.76	0.45
1:D:542:TRP:CZ3	1:D:590:MET:HB3	2.51	0.45
1:B:92:THR:HA	1:B:93:HIS:HA	1.77	0.45
1:D:140:SER:HA	1:D:450:VAL:CG2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:355:ARG:HH22	1:D:511:ARG:HH21	1.65	0.45
1:D:601:ALA:HA	1:D:653:TRP:CH2	2.52	0.45
1:B:63:TYR:CE1	1:B:172:ARG:HG2	2.52	0.45
1:D:429:ILE:O	1:D:470:CYS:HB2	2.17	0.45
1:A:412:SER:HA	1:A:413:HIS:HA	1.71	0.45
1:C:497:ARG:NE	2:C:709:HOH:O	2.34	0.45
1:D:442:TRP:CE2	1:D:488:ARG:HG2	2.52	0.45
1:C:207:THR:HG21	1:C:508:PRO:HB2	1.99	0.44
1:D:465:ASP:HB3	1:D:471:ILE:CD1	2.47	0.44
1:B:355:ARG:HH22	1:B:511:ARG:HH21	1.58	0.44
1:C:538:ARG:NH2	1:C:559:GLU:OE2	2.42	0.44
1:D:153:MET:HB2	1:D:153:MET:HE3	1.43	0.44
1:A:79:VAL:HG12	1:A:84:MET:CE	2.48	0.44
1:D:667:LYS:O	2:D:705:HOH:O	2.21	0.44
1:B:200:TYR:CZ	1:B:233:GLY:HA3	2.53	0.44
1:B:521:ASP:OD2	2:B:705:HOH:O	2.20	0.44
1:D:94:ALA:HB3	1:D:134:ASP:HB3	2.00	0.44
1:B:218:ALA:HB2	1:B:369:LYS:HE2	2.00	0.44
1:C:75:PRO:HD2	1:C:78:LEU:HD13	2.00	0.44
1:B:78:LEU:HD22	1:D:78:LEU:HD21	1.99	0.43
1:C:16:ASP:HB3	1:C:46:VAL:O	2.18	0.43
1:A:631:ILE:HG22	1:D:631:ILE:HG22	2.01	0.43
1:A:19:TRP:CD1	1:A:46:VAL:HG13	2.54	0.43
1:B:77:ARG:O	1:B:78:LEU:HD23	2.18	0.43
1:B:413:HIS:O	1:B:436:PRO:HA	2.18	0.43
1:A:85:ILE:O	1:A:192:LEU:HD12	2.19	0.43
1:B:194:THR:HG21	2:D:744:HOH:O	2.17	0.43
1:C:93:HIS:O	1:C:108:PHE:O	2.37	0.43
1:B:173:GLN:HB3	2:B:776:HOH:O	2.18	0.43
1:C:122:LEU:HD23	1:C:127:ASN:OD1	2.19	0.43
1:D:440:MET:HE1	1:D:495:TYR:OH	2.18	0.43
1:B:579:THR:OG1	1:B:596:GLN:HB2	2.18	0.43
1:A:342:SER:N	2:A:750:HOH:O	2.51	0.43
1:B:77:ARG:HG3	1:B:78:LEU:CD2	2.48	0.43
1:A:8:ASP:HA	1:C:75:PRO:HA	2.01	0.43
1:B:429:ILE:O	1:B:470:CYS:HB2	2.19	0.43
1:D:644:ASP:HB2	1:D:646:GLN:OE1	2.18	0.43
1:D:37:PRO:HA	1:D:69:GLN:OE1	2.18	0.42
1:C:426:ARG:HD3	2:C:1169:HOH:O	2.19	0.42
1:B:507:ASP:OD2	1:B:511:ARG:HD2	2.20	0.42
1:C:54:ASN:C	2:C:718:HOH:O	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:173:GLN:HB3	2:D:813:HOH:O	2.19	0.42
1:A:330:ALA:O	1:A:347:ALA:HA	2.19	0.42
1:B:316:LYS:HB2	1:B:324:TYR:CD1	2.55	0.42
1:D:318:TRP:CH2	1:D:470:CYS:HA	2.55	0.42
1:D:418:GLU:CG	1:D:464:ARG:HD2	2.50	0.42
1:D:11:ILE:H	1:D:194:THR:HB	1.83	0.42
1:D:97:VAL:HG12	1:D:104:LEU:HD12	2.01	0.42
1:D:431:ILE:HD12	1:D:470:CYS:SG	2.59	0.42
1:D:490:ARG:HD2	2:D:713:HOH:O	2.19	0.42
1:B:75:PRO:HB3	1:D:8:ASP:O	2.19	0.42
1:C:275:ARG:NH2	1:C:345:ILE:CD1	2.83	0.42
1:A:92:THR:HA	1:A:93:HIS:HA	1.83	0.42
1:C:479:GLU:CG	1:C:518:CYS:HB3	2.50	0.42
1:D:15:LEU:HD23	1:D:15:LEU:HA	1.89	0.42
1:D:412:SER:HA	1:D:413:HIS:HA	1.79	0.42
1:D:422:ASP:OD1	1:D:468:HIS:HE1	2.01	0.42
1:A:397:ASP:O	1:A:401:ILE:HG13	2.19	0.42
1:B:137:ILE:HD13	1:B:137:ILE:HG21	1.89	0.42
1:D:275:ARG:HA	1:D:275:ARG:HD2	1.85	0.42
1:A:313:ARG:HB3	2:A:950:HOH:O	2.20	0.41
1:C:334:PRO:HA	1:C:345:ILE:HD11	2.03	0.41
1:B:135:ASN:OD1	1:B:135:ASN:C	2.63	0.41
1:A:172:ARG:NH2	2:A:710:HOH:O	2.30	0.41
1:A:279:ILE:O	1:A:291:GLY:HA2	2.21	0.41
1:B:11:ILE:H	1:B:194:THR:HB	1.85	0.41
1:B:228:TYR:OH	2:B:707:HOH:O	2.21	0.41
1:C:150:THR:HG22	1:C:151:ALA:O	2.20	0.41
1:C:9:SER:OG	2:C:703:HOH:O	2.22	0.41
1:D:479:GLU:HB3	1:D:518:CYS:HB3	2.03	0.41
1:D:515:LEU:C	1:D:515:LEU:HD23	2.45	0.41
1:A:632:ARG:HA	1:A:632:ARG:HD3	1.83	0.41
1:C:433:ASP:OD2	1:C:465:ASP:OD2	2.39	0.41
1:A:601:ALA:HA	1:A:653:TRP:CH2	2.56	0.41
1:B:1:MET:HE3	1:B:1:MET:HB3	1.77	0.41
1:D:128:LEU:HD12	1:D:128:LEU:HA	1.83	0.41
1:D:303:THR:HG22	2:D:714:HOH:O	2.19	0.41
1:B:497:ARG:HD2	1:B:497:ARG:HA	1.92	0.40
1:C:619:TRP:HA	1:C:620:ASN:HA	1.89	0.40
1:D:118:VAL:O	1:D:119:THR:C	2.63	0.40
1:D:398:VAL:HG21	1:D:423:LEU:HG	2.03	0.40
1:D:589:GLU:OE2	2:D:706:HOH:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:THR:CG2	1:A:508:PRO:HB2	2.51	0.40
1:A:507:ASP:OD2	1:A:511:ARG:CD	2.70	0.40
1:D:413:HIS:O	1:D:436:PRO:HA	2.21	0.40
1:D:375:GLY:HA3	1:D:408:SER:O	2.21	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)
2:A:1166:HOH:O	2:C:1180:HOH:O[1_655]	2.16	0.04

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	612/670 (91%)	587 (96%)	25 (4%)	0	100	100
1	B	612/670 (91%)	587 (96%)	25 (4%)	0	100	100
1	C	612/670 (91%)	589 (96%)	23 (4%)	0	100	100
1	D	612/670 (91%)	591 (97%)	21 (3%)	0	100	100
All	All	2448/2680 (91%)	2354 (96%)	94 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	503/543 (93%)	498 (99%)	5 (1%)	68	81
1	B	503/543 (93%)	498 (99%)	5 (1%)	68	81
1	C	503/543 (93%)	499 (99%)	4 (1%)	73	85
1	D	503/543 (93%)	496 (99%)	7 (1%)	59	75
All	All	2012/2172 (93%)	1991 (99%)	21 (1%)	68	81

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ILE
1	A	46	VAL
1	A	74	VAL
1	A	194	THR
1	A	575	TYR
1	B	153	MET
1	B	194	THR
1	B	279	ILE
1	B	430	VAL
1	B	575	TYR
1	C	9	SER
1	C	229	ASP
1	C	387	ARG
1	C	485	ASP
1	D	9	SER
1	D	376	PHE
1	D	387	ARG
1	D	412	SER
1	D	515	LEU
1	D	568	LYS
1	D	575	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-1	ASN
1	A	102	GLN
1	A	173	GLN
1	A	350	GLN
1	A	565	ASN
1	B	-1	ASN

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Mol	Chain	Res	Type
1	B	173	GLN
1	C	276	GLN
1	C	350	GLN
1	D	350	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 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 [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	620/670 (92%)	-0.91	0 100 100	7, 14, 30, 64	0
1	B	620/670 (92%)	-0.76	2 (0%) 90 88	9, 18, 36, 65	0
1	C	620/670 (92%)	-0.86	0 100 100	7, 15, 31, 55	0
1	D	620/670 (92%)	-0.77	2 (0%) 90 88	7, 19, 37, 68	0
All	All	2480/2680 (92%)	-0.83	4 (0%) 91 90	7, 17, 34, 68	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	78	LEU	2.2
1	D	-1	ASN	2.1
1	B	274	GLY	2.0
1	B	-1	ASN	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 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.