



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2018 – 03:07 am GMT

PDB ID : 4JKM  
Title : Crystal Structure of Clostridium perfringens beta-glucuronidase  
Authors : Wallace, B.D.; Redinbo, M.R.  
Deposited on : 2013-03-09  
Resolution : 2.26 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

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix)	:	1.13
EDS	:	trunk31020
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk31020

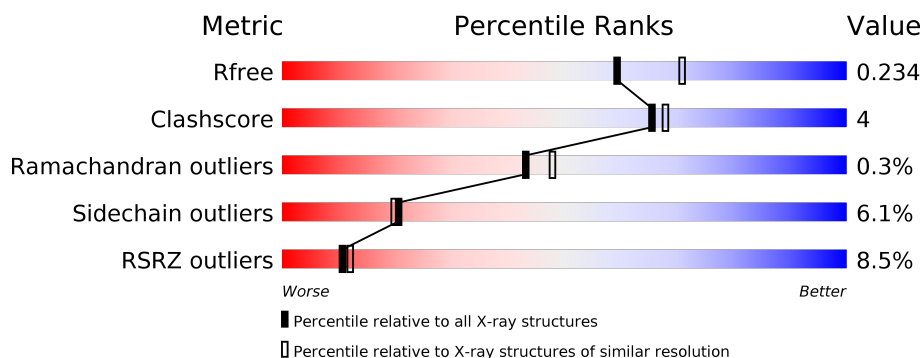
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.26 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))
$R_{free}$	111664	1178 (2.26-2.26)
Clashscore	122126	1286 (2.26-2.26)
Ramachandran outliers	120053	1253 (2.26-2.26)
Sidechain outliers	120020	1254 (2.26-2.26)
RSRZ outliers	108989	1158 (2.26-2.26)

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	602	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>13%</div> <div>.</div> </div> </div>
1	B	602	<div> <div>4%</div> <div> <div></div> <div>86%</div> <div>12%</div> <div>.</div> </div> </div>
2	C	400	<div> <div>19%</div> <div> <div></div> <div>50%</div> <div>12%</div> <div>.</div> <div>37%</div> </div> </div>
2	D	400	<div> <div>10%</div> <div> <div></div> <div>80%</div> <div>9%</div> <div>.</div> <div>10%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14718 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 Beta-glucuronidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	602	Total	C	N	O	S	0	1	0
			4863	3107	811	926	19			
1	B	602	Total	C	N	O	S	0	1	0
			4851	3098	808	926	19			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q8XP19
A	-1	ASN	-	EXPRESSION TAG	UNP Q8XP19
A	0	ALA	-	EXPRESSION TAG	UNP Q8XP19
B	-2	SER	-	EXPRESSION TAG	UNP Q8XP19
B	-1	ASN	-	EXPRESSION TAG	UNP Q8XP19
B	0	ALA	-	EXPRESSION TAG	UNP Q8XP19

- Molecule 2 is a protein called Maltose-binding periplasmic protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	362	Total	C	N	O	S	0	0	0
			2653	1693	433	521	6			
2	C	252	Total	C	N	O	S	0	0	0
			1839	1168	302	363	6			

There are 70 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MET	-	EXPRESSION TAG	UNP P0AEX9
D	2	LYS	-	EXPRESSION TAG	UNP P0AEX9
D	3	HIS	-	EXPRESSION TAG	UNP P0AEX9
D	4	HIS	-	EXPRESSION TAG	UNP P0AEX9
D	5	HIS	-	EXPRESSION TAG	UNP P0AEX9
D	6	HIS	-	EXPRESSION TAG	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
D	7	HIS	-	EXPRESSION TAG	UNP P0AEX9
D	8	HIS	-	EXPRESSION TAG	UNP P0AEX9
D	9	GLY	-	EXPRESSION TAG	UNP P0AEX9
D	25	ALA	GLY	SEE REMARK 999	UNP P0AEX9
D	376	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	377	SER	-	EXPRESSION TAG	UNP P0AEX9
D	378	SER	-	EXPRESSION TAG	UNP P0AEX9
D	379	SER	-	EXPRESSION TAG	UNP P0AEX9
D	380	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	381	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	382	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	383	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	384	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	385	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	386	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	387	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	388	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	389	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	390	ARG	-	EXPRESSION TAG	UNP P0AEX9
D	391	ASP	-	EXPRESSION TAG	UNP P0AEX9
D	392	LEU	-	EXPRESSION TAG	UNP P0AEX9
D	393	GLY	-	EXPRESSION TAG	UNP P0AEX9
D	394	THR	-	EXPRESSION TAG	UNP P0AEX9
D	395	GLU	-	EXPRESSION TAG	UNP P0AEX9
D	396	ASN	-	EXPRESSION TAG	UNP P0AEX9
D	397	LEU	-	EXPRESSION TAG	UNP P0AEX9
D	398	TYR	-	EXPRESSION TAG	UNP P0AEX9
D	399	PHE	-	EXPRESSION TAG	UNP P0AEX9
D	400	GLN	-	EXPRESSION TAG	UNP P0AEX9
C	1	MET	-	EXPRESSION TAG	UNP P0AEX9
C	2	LYS	-	EXPRESSION TAG	UNP P0AEX9
C	3	HIS	-	EXPRESSION TAG	UNP P0AEX9
C	4	HIS	-	EXPRESSION TAG	UNP P0AEX9
C	5	HIS	-	EXPRESSION TAG	UNP P0AEX9
C	6	HIS	-	EXPRESSION TAG	UNP P0AEX9
C	7	HIS	-	EXPRESSION TAG	UNP P0AEX9
C	8	HIS	-	EXPRESSION TAG	UNP P0AEX9
C	9	GLY	-	EXPRESSION TAG	UNP P0AEX9
C	25	ALA	GLY	SEE REMARK 999	UNP P0AEX9
C	376	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	377	SER	-	EXPRESSION TAG	UNP P0AEX9
C	378	SER	-	EXPRESSION TAG	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	379	SER	-	EXPRESSION TAG	UNP P0AEX9
C	380	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	381	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	382	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	383	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	384	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	385	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	386	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	387	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	388	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	389	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	390	ARG	-	EXPRESSION TAG	UNP P0AEX9
C	391	ASP	-	EXPRESSION TAG	UNP P0AEX9
C	392	LEU	-	EXPRESSION TAG	UNP P0AEX9
C	393	GLY	-	EXPRESSION TAG	UNP P0AEX9
C	394	THR	-	EXPRESSION TAG	UNP P0AEX9
C	395	GLU	-	EXPRESSION TAG	UNP P0AEX9
C	396	ASN	-	EXPRESSION TAG	UNP P0AEX9
C	397	LEU	-	EXPRESSION TAG	UNP P0AEX9
C	398	TYR	-	EXPRESSION TAG	UNP P0AEX9
C	399	PHE	-	EXPRESSION TAG	UNP P0AEX9
C	400	GLN	-	EXPRESSION TAG	UNP P0AEX9

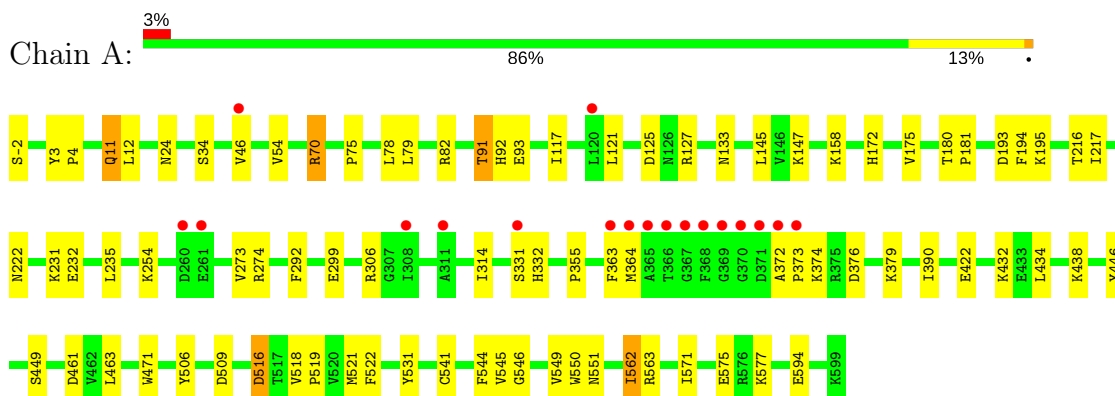
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	265	Total O 265 265	0	0
3	B	174	Total O 174 174	0	0
3	D	49	Total O 49 49	0	0
3	C	24	Total O 24 24	0	0

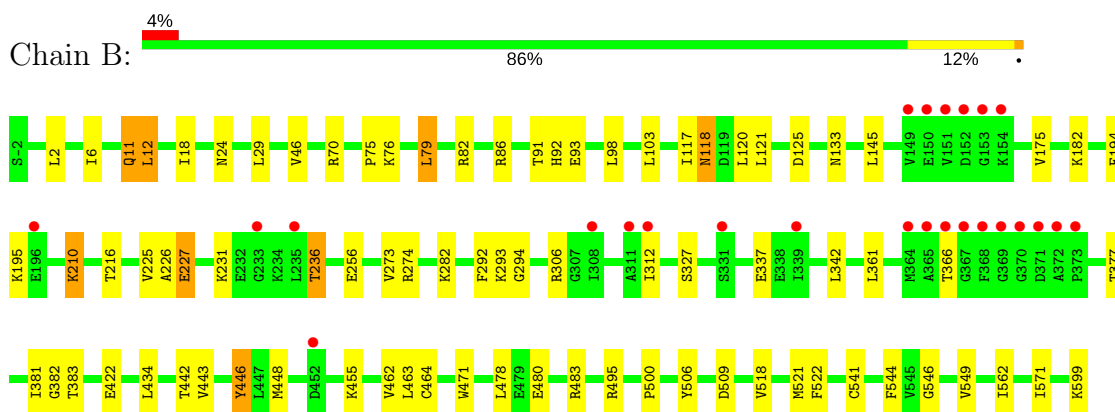
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

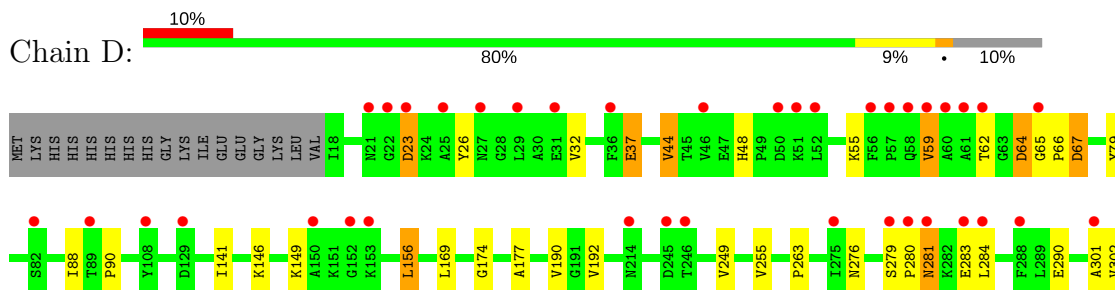
#### • Molecule 1: Beta-glucuronidase



#### • Molecule 1: Beta-glucuronidase

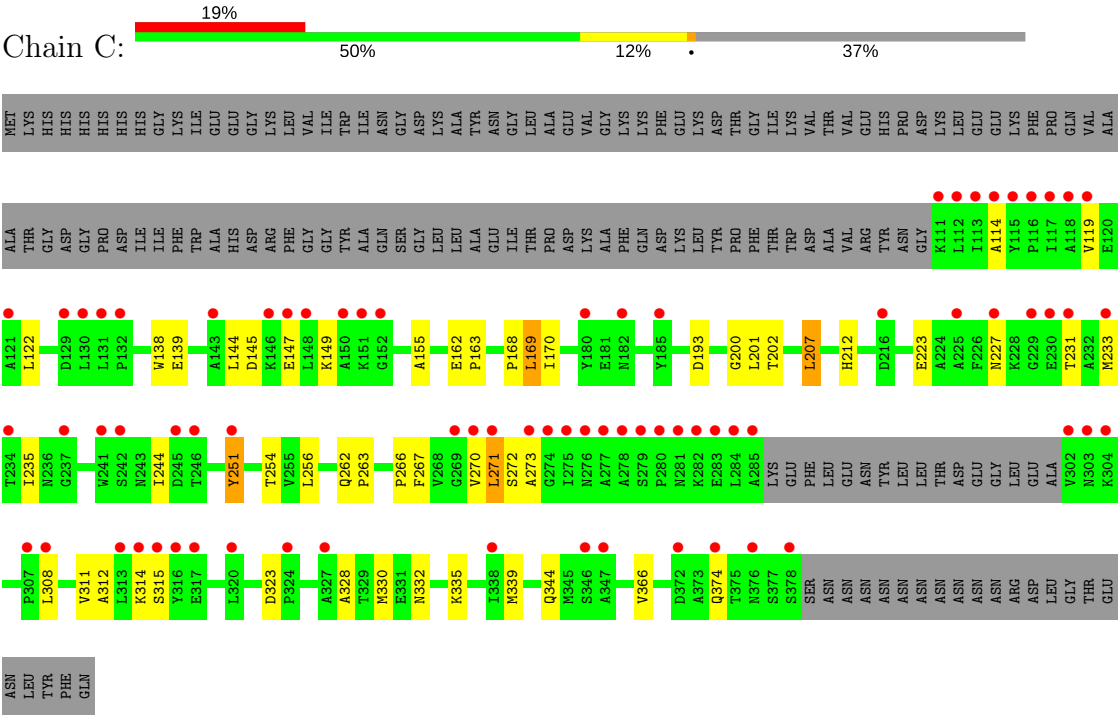


#### • Molecule 2: Maltose-binding periplasmic protein





• Molecule 2: Maltose-binding periplasmic protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.51Å 292.61Å 239.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.79 – 2.26 47.79 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.79-2.26) 99.6 (47.79-2.26)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.75 (at 2.27Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, $R_{free}$	0.207 , 0.234 0.207 , 0.234	Depositor DCC
$R_{free}$ test set	5862 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	45.8	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 48.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	14718	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.37% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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	0.26	1/4977 (0.0%)	0.45	1/6740 (0.0%)
1	B	0.23	0/4965	0.45	0/6728
2	C	0.22	0/1879	0.42	0/2569
2	D	0.23	0/2717	0.44	0/3718
All	All	0.24	1/14538 (0.0%)	0.45	1/19755 (0.0%)

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
2	D	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	373	PRO	N-CD	5.32	1.55	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	372	ALA	C-N-CD	5.42	139.77	128.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	59	VAL	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	4863	0	4766	41	0
1	B	4851	0	4733	44	0
2	C	1839	0	1730	27	0
2	D	2653	0	2441	19	0
3	A	265	0	0	4	0
3	B	174	0	0	4	0
3	C	24	0	0	3	0
3	D	49	0	0	1	0
All	All	14718	0	13670	123	0

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 (123) 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:361:LEU:H	1:B:377:THR:HG22	1.57	0.70
2:D:280:PRO:HD2	2:D:283:GLU:HB3	1.76	0.67
2:D:66:PRO:O	2:D:67:ASP:HB2	1.94	0.66
2:D:59:VAL:HG21	2:D:302:VAL:O	1.96	0.66
1:B:216:THR:HG22	1:B:227:GLU:HG3	1.78	0.65
1:B:292:PHE:HB2	1:B:546:GLY:HA3	1.77	0.64
1:B:2:LEU:O	1:B:86:ARG:NH2	2.32	0.63
1:A:292:PHE:HB2	1:A:546:GLY:HA3	1.81	0.63
2:C:223:GLU:O	2:C:227:ASN:ND2	2.30	0.62
2:D:88:ILE:HG13	2:D:90:PRO:HD3	1.82	0.62
2:D:276:ASN:HB3	2:D:280:PRO:HG3	1.82	0.62
1:A:563:ARG:NH2	3:A:810:HOH:O	2.34	0.61
2:C:308:LEU:HA	2:C:312:ALA:HA	1.85	0.58
1:A:34:SER:O	1:A:127:ARG:NH2	2.31	0.57
1:B:462:VAL:HG22	1:B:500:PRO:HG2	1.86	0.57
1:B:442:THR:HG21	1:B:464:CYS:SG	2.45	0.56
2:D:67:ASP:OD2	2:D:280:PRO:HB2	2.05	0.56
2:C:119:VAL:H	2:C:315:SER:H	1.55	0.55
2:D:141:ILE:HD13	2:D:156:LEU:HD12	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:193:ASP:HB2	2:C:374:GLN:HB2	1.90	0.54
2:D:281:ASN:ND2	3:D:533:HOH:O	2.40	0.53
1:A:516:ASP:OD1	3:A:652:HOH:O	2.19	0.53
2:C:328:ALA:O	2:C:332:ASN:ND2	2.40	0.53
2:C:114:ALA:HB3	2:C:273:ALA:HB2	1.90	0.53
1:B:337:GLU:OE1	3:B:760:HOH:O	2.19	0.53
2:D:301:ALA:HA	2:D:304:LYS:HD3	1.90	0.53
2:D:263:PRO:HB3	2:D:335:LYS:HD3	1.91	0.52
2:C:201:LEU:HD23	2:C:366:VAL:HG13	1.91	0.52
1:A:93:GLU:HB3	1:A:133:ASN:HB3	1.92	0.51
1:A:11[A]:GLN:NE2	1:B:11[A]:GLN:OE1	2.44	0.51
2:D:23:ASP:O	2:D:26:TYR:CB	2.60	0.50
1:A:575:GLU:HG2	1:A:577:LYS:HG2	1.92	0.50
1:B:93:GLU:HB3	1:B:133:ASN:HB3	1.94	0.49
1:B:480:GLU:OE2	1:B:483:ARG:NH2	2.33	0.49
2:C:235:ILE:HG23	3:C:520:HOH:O	2.12	0.49
2:D:37:GLU:CB	2:D:64:ASP:HA	2.42	0.49
2:C:263:PRO:HB3	2:C:335:LYS:HE2	1.94	0.49
1:A:332:HIS:O	1:A:355:PRO:HA	2.12	0.49
1:B:118:ASN:N	1:B:118:ASN:OD1	2.26	0.49
2:C:170:ILE:HA	2:C:200:GLY:HA3	1.94	0.49
1:B:381:ILE:HG13	1:B:383:THR:HG23	1.95	0.48
1:A:82:ARG:HH11	1:A:180:THR:HG21	1.79	0.48
1:B:562:ILE:HG12	3:B:691:HOH:O	2.13	0.48
1:A:145:LEU:HD11	1:A:147:LYS:HE3	1.96	0.48
1:A:180:THR:HG23	1:A:181:PRO:O	2.13	0.48
1:B:381:ILE:HG23	1:B:382:GLY:H	1.79	0.48
1:A:376:ASP:HB3	1:A:379:LYS:HG2	1.96	0.47
1:B:29:LEU:HD21	3:B:698:HOH:O	2.13	0.47
1:B:91:THR:HA	1:B:92:HIS:HA	1.64	0.47
2:C:254:THR:HA	2:C:328:ALA:HB1	1.97	0.47
1:A:549:VAL:HG11	1:A:571:ILE:HD11	1.96	0.47
1:A:11[B]:GLN:HG2	1:B:11[B]:GLN:NE2	2.30	0.47
1:A:217:ILE:HG13	1:A:235:LEU:HD23	1.97	0.47
1:B:361:LEU:H	1:B:377:THR:CG2	2.27	0.47
1:A:78:LEU:HD11	1:B:11[B]:GLN:HE22	1.80	0.47
1:B:471:TRP:CZ2	1:B:509:ASP:HB2	2.50	0.47
1:B:549:VAL:HG11	1:B:571:ILE:HD11	1.97	0.47
1:A:432:LYS:NZ	1:A:461:ASP:OD2	2.48	0.47
1:A:70:ARG:NH2	3:A:623:HOH:O	2.48	0.47
1:A:471:TRP:CZ2	1:A:509:ASP:HB2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:THR:HA	1:A:92:HIS:HA	1.59	0.46
1:A:216:THR:HG23	1:A:254:LYS:HB3	1.98	0.46
2:D:23:ASP:O	2:D:26:TYR:N	2.44	0.46
2:C:145:ASP:O	2:C:149:LYS:HB2	2.15	0.46
1:A:193:ASP:OD1	1:A:194:PHE:N	2.45	0.45
1:A:550:TRP:HA	1:A:551:ASN:HA	1.75	0.45
1:A:75:PRO:HG2	1:B:11[A]:GLN:HG2	1.99	0.45
2:D:32:VAL:HG21	2:D:79:TYR:OH	2.17	0.45
2:C:138:TRP:CE2	2:C:169:LEU:HG	2.52	0.45
2:C:122:LEU:HB3	3:C:520:HOH:O	2.16	0.44
1:B:256:GLU:O	3:B:685:HOH:O	2.21	0.44
2:C:207:LEU:HB2	2:C:212:HIS:HB2	1.99	0.44
2:D:65:GLY:HA3	2:D:66:PRO:HD2	1.86	0.44
2:C:271:LEU:O	2:C:273:ALA:N	2.45	0.44
2:C:202:THR:HG23	2:C:366:VAL:HG11	1.98	0.44
1:A:363:PHE:HA	1:A:364:MET:HA	1.70	0.44
1:A:562:ILE:H	1:A:562:ILE:HD13	1.82	0.44
1:B:282:LYS:NZ	1:B:599:LYS:HD2	2.33	0.44
1:B:98:LEU:HD23	1:B:120:LEU:HB2	2.00	0.44
1:B:541:CYS:HB3	1:B:544:PHE:HB2	1.99	0.44
1:B:76:LYS:HA	1:B:79:LEU:HD22	1.99	0.44
2:C:267:PHE:HB3	2:C:339:MET:HG3	2.00	0.44
2:C:244:ILE:HD11	2:C:251:TYR:CZ	2.53	0.44
2:C:271:LEU:HD12	2:C:272:SER:N	2.33	0.44
1:B:442:THR:HG22	1:B:443:VAL:N	2.33	0.43
1:A:531:TYR:OH	1:A:549:VAL:HG13	2.19	0.43
1:A:438:LYS:HE3	2:D:174:GLY:HA3	2.01	0.43
2:C:314:LYS:HA	2:C:315:SER:HA	1.53	0.43
1:B:82:ARG:HA	1:B:118:ASN:ND2	2.33	0.43
1:B:210:LYS:HG2	1:B:210:LYS:H	1.74	0.43
1:B:381:ILE:HG23	1:B:382:GLY:N	2.34	0.43
1:B:294:GLY:HA3	1:B:327:SER:O	2.19	0.43
1:B:521:MET:O	1:B:522:PHE:HB2	2.18	0.43
1:A:541:CYS:HB3	1:A:544:PHE:HB2	2.01	0.43
1:A:518:VAL:HA	1:A:519:PRO:HD3	1.89	0.42
1:B:117:ILE:O	1:B:121:LEU:HG	2.20	0.42
1:A:231:LYS:HG3	1:A:232:GLU:HG3	2.01	0.42
1:A:314:ILE:HD11	1:B:18:ILE:HD12	2.00	0.42
1:A:216:THR:HG22	3:A:723:HOH:O	2.20	0.42
1:A:521:MET:O	1:A:522:PHE:HB2	2.20	0.42
2:C:162:GLU:HA	2:C:163:PRO:HD2	1.71	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:271:LEU:HG	2:C:271:LEU:H	1.55	0.42
2:D:146:LYS:HE3	2:D:146:LYS:HB2	1.83	0.42
1:A:172:HIS:CE1	1:A:306:ARG:HB3	2.55	0.41
1:B:194:PHE:O	1:B:195:LYS:HD2	2.19	0.41
1:B:6:ILE:HG12	1:B:12:LEU:HB2	2.02	0.41
2:D:190:VAL:HG12	2:D:192:VAL:H	1.84	0.41
1:B:82:ARG:HA	1:B:118:ASN:HD21	1.85	0.41
1:A:363:PHE:CE1	1:A:364:MET:HG2	2.56	0.41
1:B:446:TYR:CE1	1:B:448:MET:HB2	2.55	0.41
2:C:168:PRO:HG3	2:C:266:PRO:HB3	2.02	0.41
1:A:299:GLU:O	1:A:306:ARG:HA	2.21	0.41
1:A:3:TYR:HA	1:A:4:PRO:HD3	1.91	0.41
1:B:293:LYS:HB2	1:B:293:LYS:HE3	1.76	0.41
2:C:155:ALA:O	2:C:233:MET:N	2.53	0.41
1:B:210:LYS:HB3	1:B:210:LYS:HE2	1.86	0.41
1:A:117:ILE:O	1:A:121:LEU:HG	2.20	0.40
1:A:11[B]:GLN:NE2	1:B:75:PRO:HG2	2.35	0.40
1:A:78:LEU:HD21	1:B:11[B]:GLN:HE22	1.86	0.40
2:C:256:LEU:HD21	3:C:520:HOH:O	2.21	0.40
2:D:59:VAL:HG22	2:D:302:VAL:HG13	2.03	0.40
1:B:226:ALA:HB1	1:B:236:THR:HG22	2.04	0.40
2:C:262:GLN:HA	2:C:263:PRO:HD3	1.91	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	601/602 (100%)	573 (95%)	28 (5%)	0	100	100
1	B	601/602 (100%)	568 (94%)	32 (5%)	1 (0%)	49	57
2	C	248/400 (62%)	227 (92%)	20 (8%)	1 (0%)	36	39

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	360/400 (90%)	338 (94%)	18 (5%)	4 (1%)	16	11
All	All	1810/2004 (90%)	1706 (94%)	98 (5%)	6 (0%)	43	48

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	270	VAL
2	D	37	GLU
2	D	44	VAL
2	D	67	ASP
2	D	177	ALA
1	B	366	THR

### 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	529/531 (100%)	499 (94%)	30 (6%)	23	23
1	B	526/531 (99%)	494 (94%)	32 (6%)	20	20
2	C	179/325 (55%)	167 (93%)	12 (7%)	18	16
2	D	251/325 (77%)	233 (93%)	18 (7%)	16	14
All	All	1485/1712 (87%)	1393 (94%)	92 (6%)	20	19

All (92) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-2	SER
1	A	11[A]	GLN
1	A	11[B]	GLN
1	A	12	LEU
1	A	24	ASN
1	A	46	VAL
1	A	54	VAL
1	A	70	ARG

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Mol	Chain	Res	Type
1	A	79	LEU
1	A	91	THR
1	A	125	ASP
1	A	158	LYS
1	A	175	VAL
1	A	195	LYS
1	A	222	ASN
1	A	273	VAL
1	A	274	ARG
1	A	331	SER
1	A	374	LYS
1	A	390	ILE
1	A	422	GLU
1	A	434	LEU
1	A	446	TYR
1	A	449	SER
1	A	463	LEU
1	A	506	TYR
1	A	516	ASP
1	A	545	VAL
1	A	562	ILE
1	A	594	GLU
1	B	11[A]	GLN
1	B	11[B]	GLN
1	B	12	LEU
1	B	24	ASN
1	B	46	VAL
1	B	70	ARG
1	B	79	LEU
1	B	103	LEU
1	B	118	ASN
1	B	125	ASP
1	B	145	LEU
1	B	175	VAL
1	B	182	LYS
1	B	210	LYS
1	B	225	VAL
1	B	227	GLU
1	B	231	LYS
1	B	236	THR
1	B	273	VAL
1	B	274	ARG

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Mol	Chain	Res	Type
1	B	306	ARG
1	B	312	ILE
1	B	342	LEU
1	B	422	GLU
1	B	434	LEU
1	B	446	TYR
1	B	455	LYS
1	B	463	LEU
1	B	478	LEU
1	B	495	ARG
1	B	506	TYR
1	B	518	VAL
2	D	23	ASP
2	D	44	VAL
2	D	48	HIS
2	D	55	LYS
2	D	62	THR
2	D	64	ASP
2	D	149	LYS
2	D	156	LEU
2	D	169	LEU
2	D	249	VAL
2	D	255	VAL
2	D	279	SER
2	D	281	ASN
2	D	284	LEU
2	D	290	GLU
2	D	325	ARG
2	D	344	GLN
2	D	374	GLN
2	C	139	GLU
2	C	144	LEU
2	C	147	GLU
2	C	169	LEU
2	C	207	LEU
2	C	231	THR
2	C	251	TYR
2	C	271	LEU
2	C	311	VAL
2	C	323	ASP
2	C	330	MET
2	C	344	GLN



Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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, 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	602/602 (100%)	0.00	18 (2%)	50	55	18, 29, 55, 162	0
1	B	602/602 (100%)	0.16	25 (4%)	36	39	24, 43, 79, 171	0
2	C	252/400 (63%)	1.41	74 (29%)	0	0	56, 82, 119, 139	0
2	D	362/400 (90%)	0.43	38 (10%)	6	6	28, 61, 104, 117	0
All	All	1818/2004 (90%)	0.33	155 (8%)	11	12	18, 44, 105, 171	0

All (155) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	368	PHE	13.5
1	B	369	GLY	13.5
1	A	371	ASP	11.0
2	C	280	PRO	11.0
1	A	365	ALA	10.2
1	B	367	GLY	10.2
1	A	367	GLY	9.5
1	B	371	ASP	9.4
2	C	273	ALA	9.1
1	A	366	THR	8.8
1	A	368	PHE	8.8
2	D	22	GLY	8.4
2	C	279	SER	8.4
1	B	365	ALA	8.0
2	D	46	VAL	7.7
1	A	364	MET	7.4
1	B	151	VAL	7.4
2	D	23	ASP	7.1
2	C	112	LEU	6.6
1	A	370	GLY	6.5
2	C	316	TYR	6.4

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Mol	Chain	Res	Type	RSRZ
2	C	111	LYS	6.2
1	B	372	ALA	6.2
2	D	284	LEU	6.1
2	C	281	ASN	6.0
1	B	370	GLY	5.9
2	D	57	PRO	5.7
2	C	132	PRO	5.7
1	A	372	ALA	5.7
2	C	302	VAL	5.7
2	C	275	ILE	5.7
2	D	27	ASN	5.6
1	B	364	MET	5.5
2	C	284	LEU	5.3
2	C	116	PRO	5.3
2	C	182	ASN	5.1
1	A	369	GLY	4.7
2	D	52	LEU	4.6
2	D	50	ASP	4.4
2	C	317	GLU	4.4
2	C	282	LYS	4.3
2	D	62	THR	4.3
2	D	31	GLU	4.2
2	D	153	LYS	4.0
2	C	278	ALA	3.8
2	D	60	ALA	3.8
1	A	363	PHE	3.8
2	C	115	TYR	3.8
1	B	235	LEU	3.7
2	C	117	ILE	3.7
1	B	366	THR	3.6
2	C	313	LEU	3.6
2	C	237	GLY	3.6
2	C	152	GLY	3.5
2	C	113	ILE	3.5
2	C	251	TYR	3.5
2	C	114	ALA	3.4
2	D	51	LYS	3.4
2	C	283	GLU	3.4
2	C	304	LYS	3.4
1	A	260	ASP	3.4
2	D	59	VAL	3.3
2	C	129	ASP	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	152	ASP	3.3
2	C	324	PRO	3.3
1	B	154	LYS	3.3
2	C	314	LYS	3.3
2	C	271	LEU	3.2
2	C	308	LEU	3.2
1	A	373	PRO	3.2
1	B	233	GLY	3.2
2	C	269	GLY	3.2
2	C	119	VAL	3.2
2	C	274	GLY	3.2
2	D	288	PHE	3.1
2	C	315	SER	3.1
2	D	245	ASP	3.1
2	C	231	THR	3.1
1	B	149	VAL	3.0
2	D	279	SER	3.0
2	C	246	THR	3.0
1	B	308	ILE	3.0
2	D	280	PRO	3.0
2	C	130	LEU	2.9
2	C	147	GLU	2.9
2	C	270	VAL	2.9
2	D	281	ASN	2.9
1	B	150	GLU	2.9
2	C	180	TYR	2.9
1	B	196	GLU	2.9
2	D	65	GLY	2.9
2	D	89	THR	2.8
2	D	275	ILE	2.8
2	C	346	SER	2.8
2	C	242	SER	2.8
1	B	311	ALA	2.8
2	D	82	SER	2.7
2	D	58	GLN	2.7
2	D	61	ALA	2.7
2	C	131	LEU	2.6
2	C	320	LEU	2.6
2	C	146	LYS	2.6
2	C	241	TRP	2.6
2	C	327	ALA	2.6
2	C	376	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	308	ILE	2.6
2	D	150	ALA	2.6
2	D	21	ASN	2.6
2	C	338	ILE	2.6
2	D	36	PHE	2.5
2	D	152	GLY	2.5
1	A	311	ALA	2.5
2	C	216	ASP	2.5
1	B	331	SER	2.5
2	C	150	ALA	2.5
2	C	277	ALA	2.4
2	C	118	ALA	2.4
1	A	261	GLU	2.4
1	B	339	ILE	2.4
2	D	214	ASN	2.4
2	C	347	ALA	2.4
2	C	307	PRO	2.3
2	C	121	ALA	2.3
2	D	25	ALA	2.3
2	D	301	ALA	2.3
2	C	148	LEU	2.3
1	B	153	GLY	2.3
2	D	246	THR	2.3
2	D	56	PHE	2.3
1	B	373	PRO	2.2
2	C	230	GLU	2.2
1	A	120	LEU	2.2
2	C	229	GLY	2.2
2	C	378	SER	2.2
2	C	303	ASN	2.1
2	D	129	ASP	2.1
2	C	276	ASN	2.1
2	D	108	TYR	2.1
2	D	283	GLU	2.1
2	C	185	TYR	2.1
1	A	331	SER	2.1
2	C	151	LYS	2.1
2	C	285	ALA	2.1
1	B	312	ILE	2.1
1	B	452	ASP	2.0
2	C	245	ASP	2.0
2	D	29	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
2	C	227	ASN	2.0
2	C	233	MET	2.0
2	C	234	THR	2.0
2	C	143	ALA	2.0
2	C	372	ASP	2.0
1	A	46	VAL	2.0
2	C	374	GLN	2.0
2	C	225	ALA	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.