



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 04:07 pm BST

PDB ID : 1OD0
Title : Family 1 b-glucosidase from *Thermotoga maritima*
Authors : Gloster, T.; Zechel, D.L.; Boraston, A.B.; Boraston, C.M.; Macdonald, J.M.;
Tilbrook, D.M.; Stick, R.V.; Davies, G.J.
Deposited on : 2003-02-12
Resolution : 2.11 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

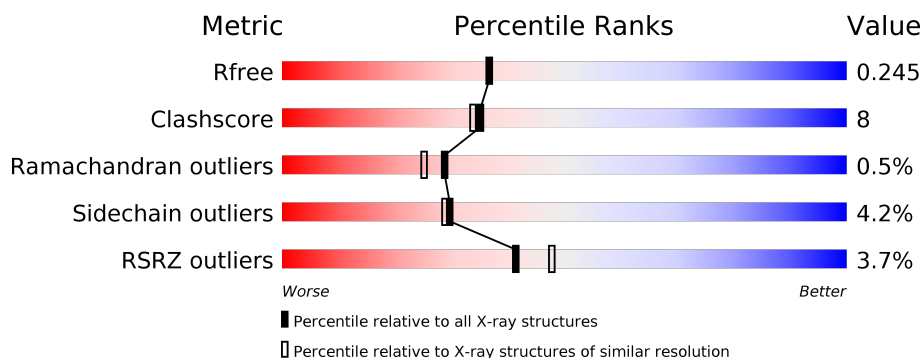
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 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	6241 (2.14-2.10)
Clashscore	141614	6778 (2.14-2.10)
Ramachandran outliers	138981	6705 (2.14-2.10)
Sidechain outliers	138945	6706 (2.14-2.10)
RSRZ outliers	127900	6112 (2.14-2.10)

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	468	<div> <div>3%</div> <div> <div></div> <div>77%</div> <div>14%</div> <div>• 7%</div> </div> </div>
1	B	468	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>17%</div> <div>• 5%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7688 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-GLUCOSIDASE A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	435	Total	C	N	O	S	0	1	0
			3550	2315	591	638	6			
1	B	446	Total	C	N	O	S	0	1	0
			3597	2340	603	647	7			

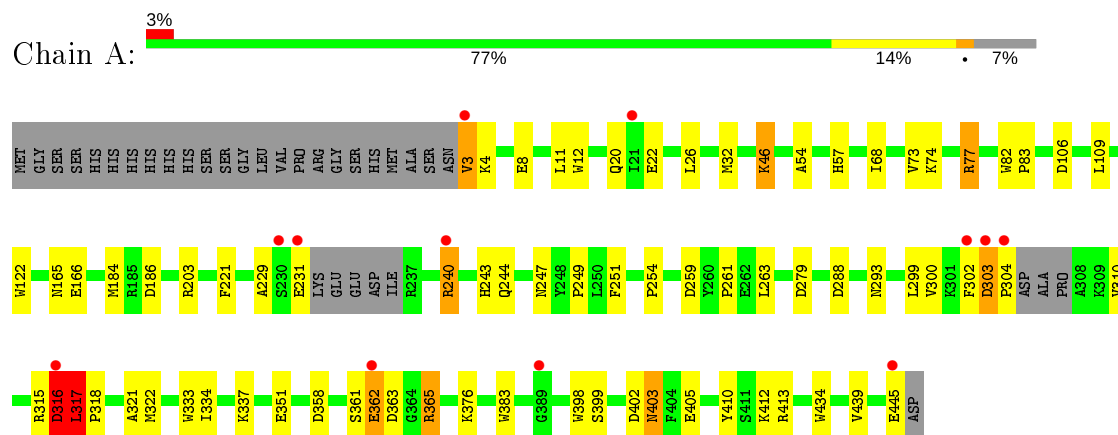
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	290	Total	O	0	0
			290	290		
2	B	251	Total	O	0	0
			251	251		

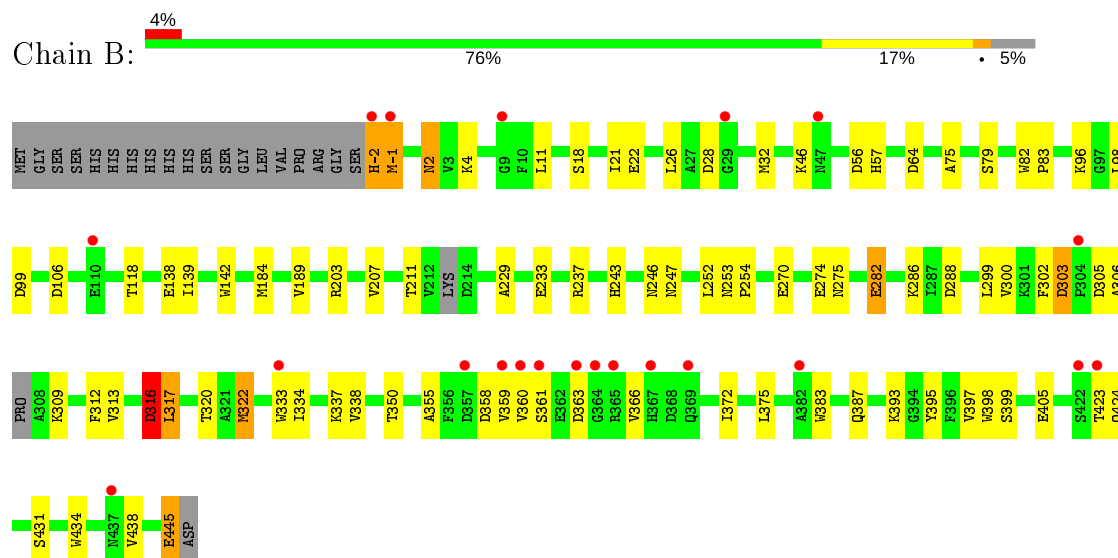
3 Residue-property plots [i](#)

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-GLUCOSIDASE A



• Molecule 1: BETA-GLUCOSIDASE A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	94.31Å 94.55Å 113.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.11 19.99 – 2.11	Depositor EDS
% Data completeness (in resolution range)	99.1 (20.00-2.11) 99.1 (19.99-2.11)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 2.11Å)	Xtriage
Refinement program	REFMAC 5.1.03	Depositor
R, R_{free}	0.210 , 0.272 0.194 , 0.245	Depositor DCC
R_{free} test set	2960 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 79.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7688	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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	0.51	0/3664	0.73	9/4980 (0.2%)
1	B	0.48	0/3711	0.73	9/5050 (0.2%)
All	All	0.50	0/7375	0.73	18/10030 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	99	ASP	CB-CG-OD2	6.79	124.41	118.30
1	A	288	ASP	CB-CG-OD2	6.42	124.08	118.30
1	A	303	ASP	CB-CG-OD2	6.07	123.76	118.30
1	B	106	ASP	CB-CG-OD2	6.07	123.76	118.30
1	A	358	ASP	CB-CG-OD2	6.05	123.75	118.30
1	B	64	ASP	CB-CG-OD2	6.02	123.72	118.30
1	B	288	ASP	CB-CG-OD2	5.97	123.67	118.30
1	A	259	ASP	CB-CG-OD2	5.96	123.66	118.30
1	A	279	ASP	CB-CG-OD2	5.83	123.55	118.30
1	A	106	ASP	CB-CG-OD2	5.76	123.48	118.30
1	B	28	ASP	CB-CG-OD2	5.72	123.45	118.30
1	B	316	ASP	CB-CG-OD2	5.68	123.41	118.30
1	B	56	ASP	CB-CG-OD2	5.65	123.39	118.30
1	A	77	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	A	186	ASP	CB-CG-OD2	5.25	123.03	118.30
1	A	316	ASP	CB-CG-OD2	5.25	123.03	118.30
1	B	303	ASP	CB-CG-OD2	5.10	122.89	118.30
1	B	305	ASP	CB-CG-OD2	5.07	122.86	118.30

There are no chirality outliers.

There are no planarity outliers.

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	3550	0	3384	50	0
1	B	3597	0	3383	64	0
2	A	290	0	0	6	0
2	B	251	0	0	8	0
All	All	7688	0	6767	114	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 8.

All (114) 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:299:LEU:HD23	1:B:313:VAL:HG22	1.36	1.05
1:B:229:ALA:HA	1:B:299:LEU:HD11	1.36	1.02
1:B:282:GLU:OE2	2:B:2196:HOH:O	1.88	0.90
1:B:274:GLU:HG3	2:B:2184:HOH:O	1.76	0.85
1:B:333:TRP:CZ2	1:B:337:LYS:HD3	2.17	0.80
1:B:270:GLU:HG2	2:B:2177:HOH:O	1.82	0.79
1:B:299:LEU:HD23	1:B:313:VAL:CG2	2.16	0.72
1:A:203:ARG:HD3	2:A:2172:HOH:O	1.93	0.69
1:B:300:VAL:HG12	1:B:312:PHE:CD1	2.28	0.68
1:B:387:GLN:NE2	1:B:387:GLN:HA	2.08	0.68
1:A:362:GLU:HG2	2:A:2242:HOH:O	1.94	0.65
1:A:299:LEU:HB2	1:A:315:ARG:HD3	1.78	0.65
1:A:300:VAL:HG13	1:A:310:VAL:HG13	1.78	0.64
1:A:376:LYS:HE3	2:A:2255:HOH:O	1.97	0.63
1:A:229:ALA:HA	1:A:299:LEU:HD11	1.83	0.59
1:B:303:ASP:O	1:B:309:LYS:HA	2.02	0.59
1:B:229:ALA:CA	1:B:299:LEU:HD11	2.22	0.58
1:B:270:GLU:HB2	2:B:2178:HOH:O	2.02	0.58
1:A:365:ARG:HD3	2:A:2245:HOH:O	2.03	0.58
1:A:12:TRP:CB	1:A:439:VAL:HG22	2.34	0.57
1:B:4:LYS:HG2	1:B:383:TRP:CD2	2.39	0.57
1:B:434:TRP:O	1:B:438:VAL:HG23	2.05	0.56
1:B:302:PHE:CE1	1:B:309:LYS:HG2	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:TRP:N	1:B:83:PRO:HD2	2.21	0.55
1:A:12:TRP:HB3	1:A:439:VAL:HG22	1.89	0.54
1:B:445:GLU:HG2	2:B:2005:HOH:O	2.07	0.54
1:A:361:SER:HB2	1:A:363:ASP:OD1	2.08	0.54
1:A:376:LYS:HG3	1:A:434:TRP:HZ2	1.73	0.53
1:A:300:VAL:HG13	1:A:310:VAL:CG1	2.39	0.52
1:A:251:PHE:C	1:A:254:PRO:HD2	2.31	0.51
1:B:302:PHE:HE1	1:B:309:LYS:HG2	1.75	0.51
1:B:375:LEU:HD23	1:B:395:TYR:OH	2.10	0.51
1:A:405:GLU:HG3	1:A:405:GLU:O	2.10	0.51
1:B:82:TRP:HB3	1:B:83:PRO:HD3	1.93	0.51
1:A:302:PHE:HE2	1:A:304:PRO:HG3	1.76	0.50
1:B:274:GLU:O	1:B:275:ASN:HB2	2.12	0.49
1:A:316:ASP:O	1:A:317:LEU:O	2.31	0.49
1:A:4:LYS:HD3	1:A:383:TRP:CE3	2.47	0.49
1:B:253:ASN:HB3	1:B:254:PRO:HD3	1.95	0.49
1:B:316:ASP:O	1:B:317:LEU:O	2.30	0.49
1:A:299:LEU:HB2	1:A:315:ARG:CD	2.41	0.48
1:B:300:VAL:HG12	1:B:312:PHE:HD1	1.74	0.48
1:A:3:VAL:HG13	2:A:2001:HOH:O	2.13	0.48
1:A:20:GLN:O	1:A:403:ASN:HB2	2.13	0.48
1:B:26:LEU:HD21	1:B:32:MET:HG2	1.95	0.48
1:B:372:ILE:HG12	1:B:431:SER:HA	1.95	0.48
1:B:2:ASN:H	1:B:2:ASN:HD22	1.62	0.48
1:B:361:SER:HB2	1:B:363:ASP:OD1	2.13	0.47
1:A:165:ASN:ND2	1:A:166:GLU:HG3	2.29	0.47
1:A:247:ASN:C	1:A:249:PRO:HD2	2.35	0.47
1:B:423:THR:O	1:B:424:GLN:HB2	2.14	0.47
1:A:321:ALA:O	1:A:412:LYS:HE3	2.14	0.47
1:B:333:TRP:HZ3	1:B:334:ILE:HD13	1.79	0.47
1:B:-2:HIS:CB	1:B:-1:MET:HE2	2.43	0.47
1:A:293:ASN:CG	1:A:351:GLU:HB2	2.35	0.47
1:B:-2:HIS:CE1	2:B:2001:HOH:O	2.67	0.47
1:B:306:ALA:C	1:B:309:LYS:H	2.19	0.46
1:B:203[A]:ARG:HG2	1:B:203[A]:ARG:HH11	1.80	0.46
1:B:79:SER:HB3	1:B:118:THR:HB	1.97	0.46
1:B:2:ASN:HD22	1:B:2:ASN:N	2.14	0.46
1:B:358:ASP:HB3	1:B:366:VAL:CG1	2.46	0.45
1:B:322:MET:HE2	1:B:322:MET:HB3	1.73	0.45
1:B:233:GLU:O	1:B:237:ARG:HG3	2.17	0.45
1:B:21:ILE:HD12	1:B:57:HIS:CG	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:393:LYS:N	1:B:393:LYS:HD2	2.31	0.45
1:B:11:LEU:HG	1:B:75:ALA:HB2	1.98	0.45
1:A:243:HIS:O	1:A:247:ASN:HB2	2.17	0.45
1:B:398:TRP:HA	1:B:399:SER:HA	1.77	0.45
1:A:82:TRP:HB3	1:A:83:PRO:HD3	2.00	0.44
1:B:-2:HIS:N	1:B:-2:HIS:CD2	2.86	0.44
1:B:-2:HIS:HB2	1:B:-1:MET:H	1.52	0.44
1:A:402:ASP:OD1	1:A:413:ARG:HB3	2.17	0.44
1:B:203[A]:ARG:HG2	1:B:203[A]:ARG:NH1	2.32	0.44
1:A:240:ARG:HG2	1:A:263:LEU:CD1	2.48	0.43
1:B:184:MET:SD	1:B:189:VAL:HG11	2.58	0.43
1:A:74:LYS:NZ	2:A:2090:HOH:O	2.51	0.43
1:A:398:TRP:HA	1:A:399:SER:HA	1.85	0.43
1:B:358:ASP:HB3	1:B:366:VAL:HG11	2.00	0.43
1:A:247:ASN:C	1:A:249:PRO:CD	2.87	0.43
1:B:-2:HIS:HB2	1:B:-1:MET:HE2	2.01	0.42
1:A:54:ALA:HA	1:A:410:TYR:OH	2.19	0.42
1:A:403:ASN:HD22	1:A:403:ASN:C	2.23	0.42
1:A:316:ASP:HB3	1:A:317:LEU:H	1.58	0.42
1:B:207:VAL:O	1:B:211:THR:HG23	2.19	0.42
1:B:317:LEU:N	1:B:317:LEU:HD23	2.34	0.42
1:B:405:GLU:O	1:B:405:GLU:HG3	2.19	0.42
1:A:322:MET:SD	1:A:412:LYS:HG3	2.60	0.42
1:A:26:LEU:HD21	1:A:32:MET:HG2	2.01	0.41
1:A:82:TRP:HB3	1:A:83:PRO:CD	2.50	0.41
1:B:-2:HIS:HE1	2:B:2001:HOH:O	2.03	0.41
1:B:320:THR:HB	1:B:355:ALA:O	2.21	0.41
1:B:274:GLU:O	1:B:275:ASN:CB	2.66	0.41
1:A:333:TRP:CZ3	1:A:334:ILE:HD12	2.55	0.41
1:A:77:ARG:NH1	1:A:351:GLU:HG3	2.35	0.41
1:B:243:HIS:O	1:B:247:ASN:HB2	2.21	0.41
1:B:333:TRP:CZ3	1:B:334:ILE:HD13	2.56	0.41
1:B:82:TRP:HB3	1:B:83:PRO:CD	2.51	0.41
1:A:244:GLN:OE1	1:A:263:LEU:HB2	2.21	0.41
1:B:139:ILE:HA	1:B:142:TRP:CE3	2.56	0.41
1:B:286:LYS:CG	2:B:2204:HOH:O	2.68	0.41
1:A:317:LEU:HA	1:A:318:PRO:HD2	1.93	0.41
1:B:18:SER:O	1:B:22:GLU:HG3	2.21	0.41
1:A:221:PHE:CE1	1:A:251:PHE:HB2	2.56	0.40
1:A:322:MET:HE3	1:A:322:MET:HB3	1.93	0.40
1:B:252:LEU:HD23	1:B:252:LEU:HA	1.83	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:LEU:HA	1:B:98:LEU:HD23	1.86	0.40
1:A:337:LYS:HD2	1:A:337:LYS:HA	1.87	0.40
1:A:68:ILE:HG23	1:A:73:VAL:HB	2.03	0.40
1:A:82:TRP:N	1:A:83:PRO:HD2	2.36	0.40
1:B:375:LEU:HD21	1:B:397:VAL:HG22	2.02	0.40
1:A:22:GLU:HA	1:A:57:HIS:HB3	2.04	0.40
1:A:122:TRP:CD1	1:A:122:TRP:N	2.90	0.40
1:A:315:ARG:HB3	1:A:317:LEU:HD22	2.02	0.40
1:A:46:LYS:HB2	1:A:46:LYS:HE2	1.30	0.40

There are no symmetry-related clashes.

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	430/468 (92%)	417 (97%)	11 (3%)	2 (0%)	29	25
1	B	441/468 (94%)	430 (98%)	9 (2%)	2 (0%)	29	25
All	All	871/936 (93%)	847 (97%)	20 (2%)	4 (0%)	29	25

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	316	ASP
1	B	316	ASP
1	A	317	LEU
1	B	317	LEU

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	362/399 (91%)	346 (96%)	16 (4%)	28	27
1	B	360/399 (90%)	346 (96%)	14 (4%)	32	32
All	All	722/798 (90%)	692 (96%)	30 (4%)	30	29

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	VAL
1	A	8	GLU
1	A	11	LEU
1	A	46	LYS
1	A	109	LEU
1	A	184	MET
1	A	231	GLU
1	A	240	ARG
1	A	261	PRO
1	A	303	ASP
1	A	316	ASP
1	A	317	LEU
1	A	362	GLU
1	A	365	ARG
1	A	403	ASN
1	A	445	GLU
1	B	-2	HIS
1	B	-1	MET
1	B	2	ASN
1	B	46	LYS
1	B	96	LYS
1	B	138	GLU
1	B	246	ASN
1	B	282	GLU
1	B	322	MET
1	B	338	VAL
1	B	350	THR

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Mol	Chain	Res	Type
1	B	359	VAL
1	B	360	VAL
1	B	445	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	246	ASN
1	A	381	GLN
1	A	403	ASN
1	A	437	ASN
1	B	2	ASN
1	B	246	ASN
1	B	381	GLN
1	B	387	GLN
1	B	441	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 carbohydrates 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 ⓘ

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	435/468 (92%)	-0.07	12 (2%) 53 59	19, 34, 50, 59	0
1	B	446/468 (95%)	0.05	21 (4%) 31 36	21, 35, 53, 66	1 (0%)
All	All	881/936 (94%)	-0.01	33 (3%) 41 48	19, 34, 53, 66	1 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	VAL	5.9
1	A	302	PHE	4.8
1	B	359	VAL	4.8
1	B	360	VAL	4.1
1	B	361	SER	3.7
1	B	333	TRP	3.6
1	A	231	GLU	3.5
1	A	303	ASP	3.5
1	B	-1	MET	3.4
1	B	9	GLY	3.2
1	B	422	SER	3.0
1	A	445	GLU	3.0
1	B	363	ASP	3.0
1	B	304	PRO	3.0
1	B	47	ASN	2.9
1	B	382	ALA	2.9
1	A	316	ASP	2.7
1	B	-2	HIS	2.7
1	A	230	SER	2.6
1	B	437	ASN	2.5
1	B	110	GLU	2.5
1	A	362	GLU	2.4
1	B	357	ASP	2.3
1	B	369	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	367	HIS	2.2
1	A	240	ARG	2.1
1	B	365	ARG	2.1
1	A	389	GLY	2.1
1	B	364	GLY	2.1
1	B	29	GLY	2.0
1	A	21	ILE	2.0
1	B	423	THR	2.0
1	A	304	PRO	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.