



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:48 AM UTC

PDB ID : 3WO8 / pdb_00003wo8
Title : Crystal structure of the beta-N-acetylglucosaminidase from *Thermotoga maritima*
Authors : Mine, S.; Kado, Y.; Watanabe, M.; Inoue, T.; Ishikawa, K.
Deposited on : 2013-12-20
Resolution : 2.43 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

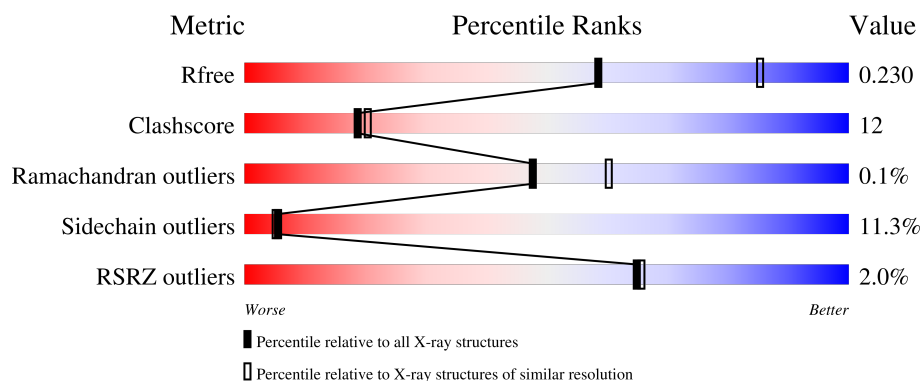
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.43 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	2340 (2.46-2.42)
Clashscore	190562	2400 (2.46-2.42)
Ramachandran outliers	187476	2379 (2.46-2.42)
Sidechain outliers	187428	2379 (2.46-2.42)
RSRZ outliers	180081	2340 (2.46-2.42)

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	478	
1	B	478	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7196 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-N-acetylglucosaminidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	447	Total	C	N	O	S	0	0	0
			3540	2285	575	668	12			
1	B	448	Total	C	N	O	S	0	0	0
			3549	2291	576	670	12			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	expression tag	UNP Q9WZR6
A	-9	ASN	-	expression tag	UNP Q9WZR6
A	-8	HIS	-	expression tag	UNP Q9WZR6
A	-7	LYS	-	expression tag	UNP Q9WZR6
A	-6	VAL	-	expression tag	UNP Q9WZR6
A	-5	HIS	-	expression tag	UNP Q9WZR6
A	-4	HIS	-	expression tag	UNP Q9WZR6
A	-3	HIS	-	expression tag	UNP Q9WZR6
A	-2	HIS	-	expression tag	UNP Q9WZR6
A	-1	HIS	-	expression tag	UNP Q9WZR6
A	0	HIS	-	expression tag	UNP Q9WZR6
B	-10	MET	-	expression tag	UNP Q9WZR6
B	-9	ASN	-	expression tag	UNP Q9WZR6
B	-8	HIS	-	expression tag	UNP Q9WZR6
B	-7	LYS	-	expression tag	UNP Q9WZR6
B	-6	VAL	-	expression tag	UNP Q9WZR6
B	-5	HIS	-	expression tag	UNP Q9WZR6
B	-4	HIS	-	expression tag	UNP Q9WZR6
B	-3	HIS	-	expression tag	UNP Q9WZR6
B	-2	HIS	-	expression tag	UNP Q9WZR6
B	-1	HIS	-	expression tag	UNP Q9WZR6
B	0	HIS	-	expression tag	UNP Q9WZR6

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	60	Total 60	O 60	0	0
2	B	47	Total 47	O 47	0	0

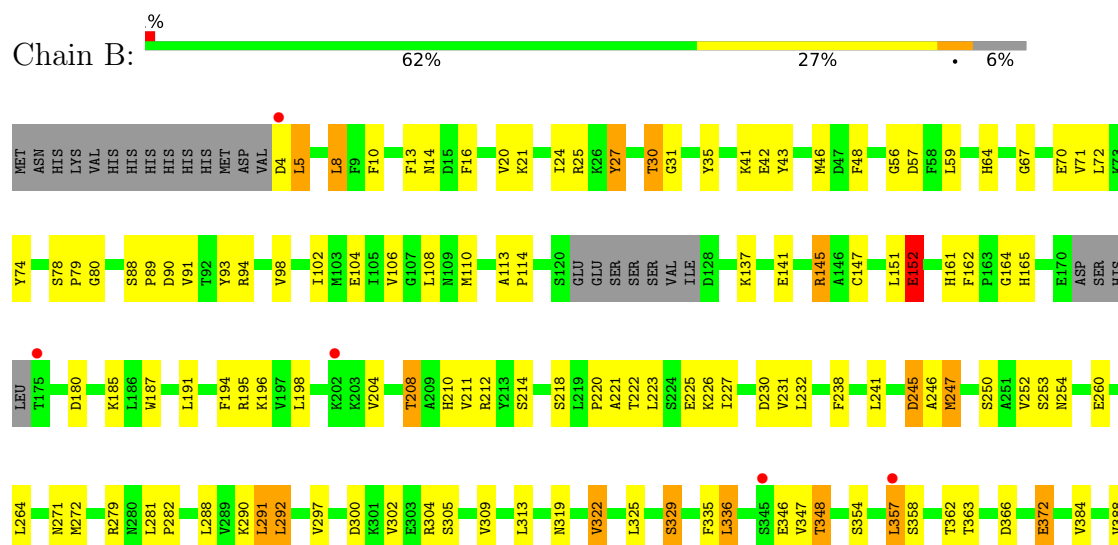
3 Residue-property plots [i](#)

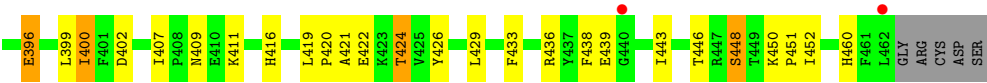
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Beta-N-acetylglucosaminidase



• Molecule 1: Beta-N-acetylglucosaminidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	133.46Å 133.46Å 142.33Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.56 – 2.43 26.56 – 2.43	Depositor EDS
% Data completeness (in resolution range)	92.0 (26.56-2.43) 92.0 (26.56-2.43)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.14 (at 2.44Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.194 , 0.246 (Not available) , 0.230	Depositor DCC
R_{free} test set	2572 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	40.4	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.036 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7196	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.12% 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	0.99	1/3614 (0.0%)	1.15	9/4881 (0.2%)
1	B	0.92	0/3623	1.11	12/4894 (0.2%)
All	All	0.95	1/7237 (0.0%)	1.13	21/9775 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	370	ILE	CA-CB	-5.93	1.51	1.54

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	450	LYS	CA-C-N	8.64	128.23	119.24
1	A	450	LYS	C-N-CA	8.64	128.23	119.24
1	B	27	TYR	N-CA-C	-7.36	99.01	110.42
1	B	78	SER	CA-C-N	-6.71	112.89	119.87
1	B	78	SER	C-N-CA	-6.71	112.89	119.87
1	B	56	GLY	N-CA-C	6.34	119.14	110.58
1	B	421	ALA	N-CA-C	6.28	120.04	112.38
1	B	27	TYR	CA-C-N	-6.14	115.72	123.15
1	B	27	TYR	C-N-CA	-6.14	115.72	123.15
1	B	245	ASP	N-CA-C	-6.09	102.92	110.41
1	A	231	VAL	N-CA-C	5.66	116.68	111.81
1	A	80	GLY	N-CA-C	5.64	117.80	112.04
1	B	152	GLU	N-CA-C	-5.59	105.19	111.28
1	B	204	VAL	N-CA-C	5.56	118.43	110.09
1	A	359	GLN	N-CA-C	5.56	117.34	111.28
1	B	80	GLY	N-CA-C	5.50	119.55	112.18
1	B	446	THR	N-CA-C	-5.35	102.83	110.59
1	A	60	ILE	N-CA-C	5.28	115.70	107.99
1	A	304	ARG	N-CA-C	-5.14	105.57	111.07
1	A	28	ARG	CA-C-N	-5.01	114.75	119.76
1	A	28	ARG	C-N-CA	-5.01	114.75	119.76

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	3540	0	3548	78	0
1	B	3549	0	3559	95	0
2	A	60	0	0	1	0
2	B	47	0	0	3	0
All	All	7196	0	7107	169	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 12.

All (169) 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:59:LEU:HD13	1:B:110:MET:HE1	1.53	0.91
1:A:414:LYS:HA	1:A:414:LYS:HE2	1.50	0.90
1:A:59:LEU:HD13	1:A:110:MET:HE1	1.54	0.88
1:A:43:TYR:OH	1:A:372:GLU:OE2	1.97	0.82
1:B:402:ASP:OD2	1:B:416:HIS:HD2	1.67	0.78
1:A:210:HIS:CE1	1:A:245:ASP:HB3	2.20	0.76
1:B:222:THR:HG21	1:B:247:MET:HE1	1.67	0.75
1:B:5:LEU:HD23	1:B:5:LEU:H	1.51	0.75
1:B:227:ILE:O	1:B:231:VAL:HG12	1.87	0.75
1:B:64:HIS:CD2	1:B:72:LEU:H	2.04	0.74
1:B:407:ILE:HD12	1:B:407:ILE:O	1.89	0.73
1:B:247:MET:HA	1:B:247:MET:HE2	1.72	0.71
1:B:357:LEU:HB2	2:B:515:HOH:O	1.89	0.71
1:A:148:GLU:O	1:A:152:GLU:HG2	1.91	0.70
1:A:347:VAL:O	1:A:380:VAL:O	2.08	0.70
1:A:400:ILE:HB	1:A:424:THR:HG22	1.73	0.70
1:A:46:MET:HG2	1:A:322:VAL:HG13	1.74	0.70
1:A:226:LYS:O	1:A:230:ASP:HB2	1.92	0.69
1:A:64:HIS:HD2	1:A:72:LEU:H	1.40	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:LEU:HD12	1:B:271:ASN:O	1.94	0.68
1:B:57:ASP:OD2	1:B:313:LEU:HG	1.94	0.67
1:A:68:GLN:HG2	1:A:69:LEU:HD13	1.76	0.66
1:A:414:LYS:HE2	1:A:414:LYS:CA	2.25	0.66
1:B:64:HIS:CD2	1:B:71:VAL:H	2.13	0.66
1:B:64:HIS:HD2	1:B:72:LEU:H	1.43	0.65
1:A:247:MET:HA	1:A:247:MET:CE	2.27	0.65
1:A:208:THR:HG23	1:A:222:THR:OG1	1.97	0.65
1:B:419:LEU:HB3	1:B:424:THR:HG21	1.78	0.65
1:B:24:ILE:O	1:B:27:TYR:O	2.15	0.64
1:B:227:ILE:HA	1:B:231:VAL:HG12	1.79	0.64
1:A:64:HIS:CD2	1:A:72:LEU:H	2.17	0.61
1:B:407:ILE:O	1:B:407:ILE:CD1	2.48	0.61
1:A:21:LYS:HG2	1:A:51:PHE:CE1	2.35	0.61
1:A:339:GLU:O	1:A:340:LYS:HG3	2.00	0.60
1:A:367:TYR:HA	1:A:370:ILE:HD12	1.83	0.60
1:A:402:ASP:OD2	1:A:416:HIS:HD2	1.85	0.60
1:B:208:THR:HG23	1:B:222:THR:OG1	2.02	0.59
1:A:184:GLU:O	1:A:188:GLU:HG2	2.03	0.59
1:B:363:THR:O	1:B:366:ASP:HB2	2.03	0.59
1:B:89:PRO:HB3	1:B:145:ARG:HG3	1.84	0.59
1:B:247:MET:HE2	1:B:252:VAL:HG21	1.84	0.58
1:B:291:LEU:HD13	1:B:297:VAL:HG23	1.84	0.58
1:B:191:LEU:O	1:B:195:ARG:HG3	2.04	0.58
1:B:211:VAL:O	1:B:220:PRO:HA	2.03	0.57
1:B:247:MET:HA	1:B:247:MET:CE	2.35	0.57
1:A:247:MET:HA	1:A:247:MET:HE2	1.85	0.57
1:B:438:PHE:O	1:B:443:ILE:HD11	2.04	0.57
1:A:359:GLN:HE22	1:B:14:ASN:HD21	1.49	0.57
1:B:210:HIS:HD2	1:B:222:THR:OG1	1.88	0.57
1:A:117:ASP:OD2	1:A:161:HIS:HD2	1.88	0.57
1:B:420:PRO:O	1:B:424:THR:HG23	2.04	0.57
1:B:43:TYR:OH	1:B:372:GLU:OE2	2.22	0.57
1:B:88:SER:O	1:B:91:VAL:HG22	2.05	0.56
1:B:433:PHE:CZ	1:B:436:ARG:HD3	2.40	0.56
1:A:348:THR:HG22	1:A:382:ASN:HB2	1.88	0.56
1:B:14:ASN:HD22	1:B:35:TYR:HB2	1.72	0.55
1:A:210:HIS:HD2	1:A:222:THR:OG1	1.88	0.55
1:B:416:HIS:HE1	2:B:509:HOH:O	1.90	0.55
1:A:414:LYS:HA	1:A:414:LYS:CE	2.29	0.54
1:A:347:VAL:HG12	1:A:399:LEU:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:TYR:HA	1:B:74:TYR:HA	1.91	0.53
1:B:433:PHE:CE1	1:B:436:ARG:CD	2.92	0.53
1:A:164:GLY:O	1:A:213:TYR:OH	2.21	0.53
1:A:167:LYS:HE2	1:A:190:ASP:OD1	2.09	0.52
1:B:180:ASP:HA	1:B:214:SER:OG	2.09	0.52
1:B:220:PRO:HG2	1:B:223:LEU:HD12	1.90	0.52
1:B:5:LEU:HD12	1:B:288:LEU:HD23	1.92	0.52
1:A:227:ILE:HA	1:A:231:VAL:HG13	1.92	0.52
1:A:200:ARG:HH21	1:A:200:ARG:HG3	1.74	0.52
1:A:419:LEU:HB3	1:A:424:THR:HG21	1.92	0.51
1:B:90:ASP:OD1	1:B:145:ARG:NH1	2.44	0.51
1:B:198:LEU:HD22	1:B:238:PHE:CD1	2.45	0.51
1:B:433:PHE:CE1	1:B:436:ARG:NE	2.79	0.51
1:A:59:LEU:HD13	1:A:110:MET:CE	2.34	0.51
1:A:64:HIS:CD2	1:A:71:VAL:H	2.29	0.51
1:B:260:GLU:O	1:B:264:LEU:HG	2.11	0.51
1:A:233:ARG:O	1:A:237:GLY:HA2	2.11	0.51
1:A:198:LEU:HD22	1:A:238:PHE:CD1	2.47	0.50
1:A:345:SER:O	1:A:379:GLU:HB2	2.11	0.50
1:B:402:ASP:OD2	1:B:416:HIS:CD2	2.57	0.50
1:A:346:GLU:HA	1:A:379:GLU:O	2.10	0.50
1:A:187:TRP:CH2	1:A:195:ARG:HD3	2.46	0.49
1:B:250:SER:HB2	1:B:254:ASN:ND2	2.27	0.49
1:A:25:ARG:HG2	1:A:55:GLU:OE1	2.13	0.49
1:A:210:HIS:HE1	1:A:246:ALA:O	1.94	0.49
1:A:344:THR:HG21	1:A:398:ASP:HB3	1.95	0.48
1:B:113:ALA:HB1	1:B:114:PRO:HA	1.95	0.48
1:B:291:LEU:CD1	1:B:297:VAL:HG23	2.44	0.48
1:B:161:HIS:O	1:B:164:GLY:N	2.46	0.48
1:A:343:HIS:NE2	1:A:377:PHE:O	2.46	0.48
1:B:46:MET:HG2	1:B:322:VAL:HG13	1.94	0.48
1:B:30:THR:HG22	1:B:272:MET:HE1	1.95	0.48
1:A:79:PRO:O	1:B:448:SER:OG	2.30	0.47
1:B:227:ILE:CA	1:B:231:VAL:HG12	2.44	0.47
1:B:325:LEU:O	1:B:329:SER:OG	2.32	0.47
1:A:247:MET:HA	1:A:247:MET:HE3	1.95	0.47
1:B:16:PHE:CD1	1:B:48:PHE:CD1	3.03	0.47
1:B:210:HIS:HE1	1:B:246:ALA:O	1.97	0.47
1:A:210:HIS:CD2	1:A:222:THR:OG1	2.68	0.47
1:A:380:VAL:O	1:A:382:ASN:N	2.42	0.47
1:B:64:HIS:HD2	1:B:72:LEU:N	2.09	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:HIS:CD2	1:A:72:LEU:HB2	2.49	0.47
1:A:247:MET:HE2	1:A:252:VAL:HG21	1.95	0.47
1:A:341:ILE:HG22	1:A:342:ASP:N	2.29	0.47
1:B:362:THR:OG1	1:B:450:LYS:NZ	2.48	0.46
1:B:102:ILE:O	1:B:106:VAL:HG23	2.15	0.46
1:B:227:ILE:C	1:B:231:VAL:HG12	2.40	0.46
1:B:59:LEU:HD13	1:B:110:MET:CE	2.35	0.46
1:B:10:PHE:CG	1:B:31:GLY:HA3	2.50	0.46
1:A:156:ILE:CD1	1:A:316:ALA:HB2	2.46	0.46
1:B:208:THR:HG22	1:B:245:ASP:H	1.80	0.45
1:A:135:ASP:HB3	1:A:138:ILE:HD12	1.98	0.45
1:B:433:PHE:CZ	1:B:436:ARG:CD	2.99	0.45
1:A:225:GLU:HA	1:A:268:ALA:HB1	1.97	0.45
1:B:292:LEU:HD11	1:B:302:VAL:HG21	1.98	0.45
1:A:448:SER:HB2	1:B:79:PRO:O	2.17	0.45
1:B:14:ASN:HD22	1:B:35:TYR:CB	2.30	0.45
1:B:162:PHE:CE2	1:B:232:LEU:HB2	2.52	0.45
1:A:380:VAL:O	1:A:381:GLU:HB3	2.16	0.45
1:B:279:ARG:O	1:B:282:PRO:HD2	2.16	0.45
1:A:244:SER:OG	1:A:245:ASP:O	2.34	0.44
1:A:344:THR:O	1:A:345:SER:C	2.58	0.44
1:A:421:ALA:O	1:A:442:LYS:HE3	2.17	0.44
1:B:16:PHE:CE1	1:B:48:PHE:CD1	3.06	0.44
1:B:93:TYR:OH	1:B:152:GLU:HG3	2.18	0.44
1:B:400:ILE:HB	1:B:424:THR:HB	1.99	0.44
1:A:10:PHE:CG	1:A:31:GLY:HA3	2.52	0.44
1:A:412:ALA:O	1:A:413:LEU:C	2.61	0.44
1:B:67:GLY:O	1:B:70:GLU:HG3	2.18	0.44
1:A:341:ILE:CG2	1:A:342:ASP:N	2.81	0.44
1:B:221:ALA:O	1:B:222:THR:C	2.59	0.44
1:B:194:PHE:O	1:B:198:LEU:HB2	2.18	0.43
1:B:147:CYS:O	1:B:151:LEU:HD12	2.17	0.43
1:B:208:THR:HG21	1:B:222:THR:HG23	2.00	0.43
1:B:416:HIS:O	1:B:426:TYR:OH	2.27	0.43
1:B:335:PHE:HD2	2:B:529:HOH:O	2.00	0.43
1:B:241:LEU:HD11	1:B:309:VAL:HG22	1.99	0.43
1:B:187:TRP:CH2	1:B:195:ARG:HD3	2.53	0.43
1:A:165:HIS:CD2	1:A:165:HIS:O	2.72	0.43
1:A:341:ILE:CG2	1:A:342:ASP:H	2.32	0.43
1:A:203:LYS:HE3	1:A:315:PHE:CG	2.54	0.42
1:B:41:LYS:O	1:B:42:GLU:C	2.61	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:MET:CG	1:A:322:VAL:HG13	2.47	0.42
1:B:227:ILE:O	1:B:231:VAL:CG1	2.64	0.42
1:B:336:LEU:HB2	1:B:443:ILE:HB	2.02	0.42
1:B:210:HIS:CD2	1:B:222:THR:OG1	2.71	0.42
1:A:64:HIS:HD2	1:A:71:VAL:H	1.64	0.42
1:A:263:SER:HB3	1:A:291:LEU:HD13	2.01	0.42
1:A:193:PRO:O	1:A:197:VAL:HG23	2.19	0.42
1:A:208:THR:CG2	1:A:222:THR:OG1	2.68	0.42
1:A:195:ARG:O	1:A:199:GLU:HB3	2.20	0.42
1:B:13:PHE:CD2	1:B:20:VAL:HG13	2.55	0.42
1:A:208:THR:HG21	1:A:222:THR:HG23	2.02	0.41
1:A:305:SER:O	1:A:306:ILE:C	2.63	0.41
1:B:8:LEU:HD11	1:B:305:SER:HB3	2.02	0.41
1:B:281:LEU:O	1:B:282:PRO:C	2.64	0.41
1:A:64:HIS:HE1	2:A:502:HOH:O	2.03	0.41
1:A:211:VAL:HG12	1:A:212:ARG:N	2.35	0.41
1:B:247:MET:O	1:B:252:VAL:HG23	2.20	0.41
1:B:212:ARG:CD	1:B:218:SER:O	2.68	0.41
1:B:348:THR:HG21	1:B:396:GLU:O	2.20	0.41
1:B:281:LEU:N	1:B:282:PRO:CD	2.83	0.41
1:B:226:LYS:O	1:B:230:ASP:HB2	2.21	0.41
1:A:346:GLU:OE1	1:A:381:GLU:HG2	2.21	0.40
1:B:187:TRP:CZ2	1:B:195:ARG:HD3	2.56	0.40
1:B:210:HIS:CE1	1:B:245:ASP:HB3	2.56	0.40
1:A:431:ASN:OD1	1:A:433:PHE:N	2.50	0.40
1:A:432:PRO:O	1:A:435:VAL:HG23	2.20	0.40
1:B:21:LYS:O	1:B:25:ARG:HG3	2.22	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	441/478 (92%)	418 (95%)	22 (5%)	1 (0%)	43	53
1	B	442/478 (92%)	423 (96%)	19 (4%)	0	100	100
All	All	883/956 (92%)	841 (95%)	41 (5%)	1 (0%)	48	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	410	GLU

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	393/423 (93%)	354 (90%)	39 (10%)	7	7
1	B	394/423 (93%)	344 (87%)	50 (13%)	4	3
All	All	787/846 (93%)	698 (89%)	89 (11%)	5	5

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	23	ILE
1	A	36	PRO
1	A	57	ASP
1	A	69	LEU
1	A	88	SER
1	A	91	VAL
1	A	94	ARG
1	A	98	VAL
1	A	104	GLU
1	A	137	LYS
1	A	160	LYS
1	A	186	LEU
1	A	188	GLU
1	A	201	GLU

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Mol	Chain	Res	Type
1	A	208	THR
1	A	216	ILE
1	A	225	GLU
1	A	231	VAL
1	A	247	MET
1	A	252	VAL
1	A	253	SER
1	A	259	GLU
1	A	322	VAL
1	A	336	LEU
1	A	338	PHE
1	A	339	GLU
1	A	343	HIS
1	A	345	SER
1	A	346	GLU
1	A	372	GLU
1	A	379	GLU
1	A	384	VAL
1	A	389	GLU
1	A	414	LYS
1	A	418	SER
1	A	422	GLU
1	A	442	LYS
1	A	448	SER
1	B	4	ASP
1	B	5	LEU
1	B	8	LEU
1	B	30	THR
1	B	94	ARG
1	B	98	VAL
1	B	104	GLU
1	B	108	LEU
1	B	137	LYS
1	B	141	GLU
1	B	145	ARG
1	B	152	GLU
1	B	165	HIS
1	B	185	LYS
1	B	196	LYS
1	B	208	THR
1	B	225	GLU
1	B	247	MET

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Mol	Chain	Res	Type
1	B	253	SER
1	B	290	LYS
1	B	291	LEU
1	B	292	LEU
1	B	300	ASP
1	B	304	ARG
1	B	319	ASN
1	B	322	VAL
1	B	329	SER
1	B	336	LEU
1	B	346	GLU
1	B	347	VAL
1	B	348	THR
1	B	354	SER
1	B	357	LEU
1	B	358	SER
1	B	372	GLU
1	B	384	VAL
1	B	388	VAL
1	B	396	GLU
1	B	399	LEU
1	B	400	ILE
1	B	409	ASN
1	B	411	LYS
1	B	422	GLU
1	B	424	THR
1	B	429	LEU
1	B	439	GLU
1	B	448	SER
1	B	451	PRO
1	B	452	ILE
1	B	460	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	64	HIS
1	A	142	HIS
1	A	161	HIS
1	A	165	HIS
1	A	210	HIS

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Mol	Chain	Res	Type
1	A	254	ASN
1	A	255	ASN
1	A	298	GLN
1	B	14	ASN
1	B	64	HIS
1	B	210	HIS
1	B	255	ASN
1	B	319	ASN
1	B	343	HIS
1	B	416	HIS

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	447/478 (93%)	-0.33	11 (2%) 58 59	25, 42, 89, 141	0
1	B	448/478 (93%)	-0.13	7 (1%) 70 72	28, 50, 85, 108	0
All	All	895/956 (93%)	-0.23	18 (2%) 65 66	25, 47, 87, 141	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	433	PHE	6.1
1	A	341	ILE	4.2
1	A	380	VAL	3.1
1	B	175	THR	3.0
1	A	461	PHE	3.0
1	B	462	LEU	3.0
1	A	344	THR	2.8
1	B	202	LYS	2.7
1	B	4	ASP	2.6
1	A	345	SER	2.6
1	A	381	GLU	2.5
1	B	345	SER	2.5
1	A	354	SER	2.4
1	A	346	GLU	2.3
1	A	440	GLY	2.3
1	A	175	THR	2.1
1	B	440	GLY	2.1
1	B	357	LEU	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.