



Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 07:08 PM BST

PDB ID : 2CBH
Title : DETERMINATION OF THE THREE-DIMENSIONAL STRUCTURE OF THE C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I FROM TRICHODERMA REESEI. A STUDY USING NUCLEAR MAGNETIC RESONANCE AND HYBRID DISTANCE GEOMETRY-DYNAMICAL SIMULATED ANNEALING
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Deposited on : 1989-05-30

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/NMRValidationReportHelp>
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:

Cyrange : Kirchner and Güntert (2011)
NmrClust : Kelley et al. (1996)
MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : rb-20027457
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027457

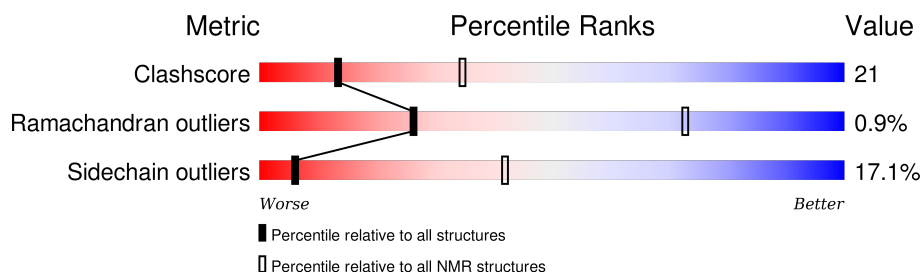
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

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)	NMR archive (#Entries)
Clashscore	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	36	

2 Ensemble composition and analysis

This entry contains 41 models. Model 32 is the overall representative, medoid model (most similar to other models).

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:2-A:36 (35)	0.21	32

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 8 clusters and 6 single-model clusters were found.

Cluster number	Models
1	2, 4, 5, 6, 22, 23, 27, 28, 29, 35, 37, 40
2	11, 19, 24, 25, 26, 31, 32, 36
3	7, 18, 20, 39
4	9, 13, 38
5	1, 15
6	3, 12
7	8, 33
8	14, 21
Single-model clusters	10; 16; 17; 30; 34; 41

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 495 atoms, of which 235 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I.

Mol	Chain	Residues	Atoms						Trace
1	A	36	Total	C	H	N	O	S	0
			495	159	235	43	54	4	

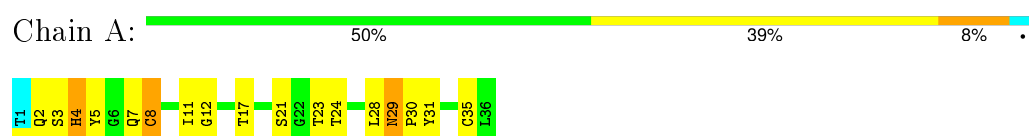
4.2.3 Score per residue for model 3

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



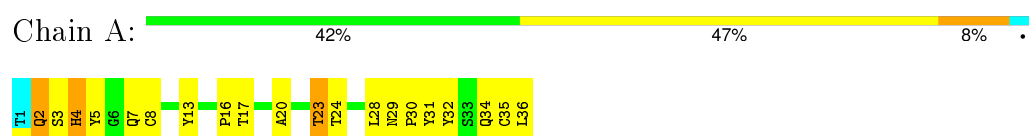
4.2.4 Score per residue for model 4

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.5 Score per residue for model 5

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.6 Score per residue for model 6

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.7 Score per residue for model 7

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



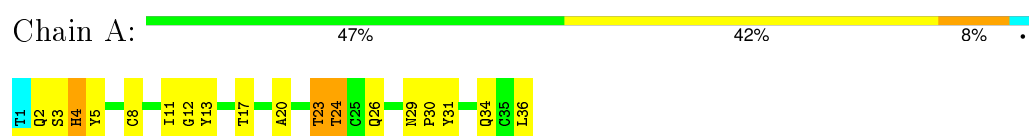
4.2.8 Score per residue for model 8

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.9 Score per residue for model 9

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



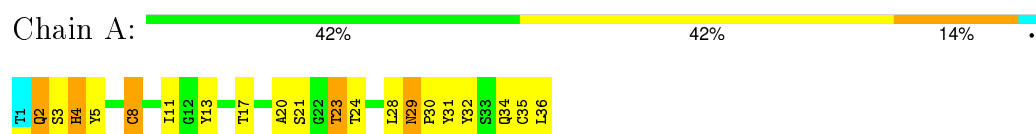
4.2.10 Score per residue for model 10

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



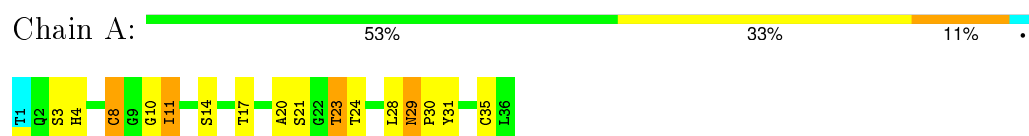
4.2.11 Score per residue for model 11

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.12 Score per residue for model 12

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.13 Score per residue for model 13

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I

Chain A:  56% 33% 8% •



4.2.14 Score per residue for model 14

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I

Chain A:  56% 31% 11% •



4.2.15 Score per residue for model 15

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I

Chain A:  58% 33% 6% •



4.2.16 Score per residue for model 16

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I

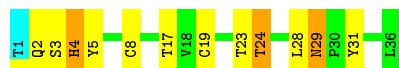
Chain A:  56% 36% 6% •



4.2.17 Score per residue for model 17

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I

Chain A:  64% 25% 8% •



4.2.18 Score per residue for model 18

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I

Chain A:  58% 33% 6% •



4.2.19 Score per residue for model 19

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I

Chain A:  58% 33% 6% •



4.2.20 Score per residue for model 20

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I

Chain A:  64% 19% 14% •



4.2.21 Score per residue for model 21

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I

Chain A:  58% 36% • •



4.2.22 Score per residue for model 22

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I

Chain A:  64% 28% 6% •



4.2.23 Score per residue for model 23

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



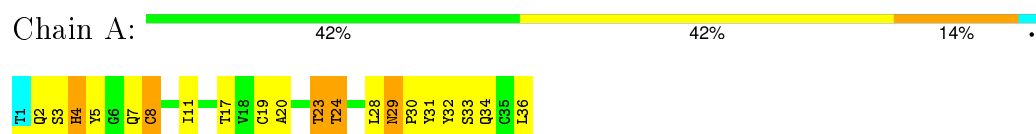
4.2.24 Score per residue for model 24

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



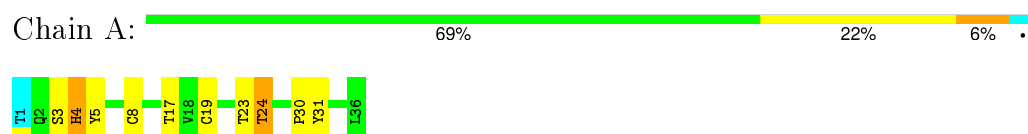
4.2.25 Score per residue for model 25

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.26 Score per residue for model 26

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.27 Score per residue for model 27

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



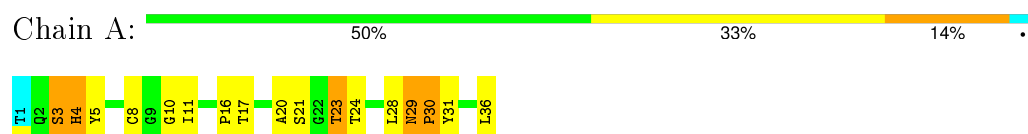
4.2.28 Score per residue for model 28

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



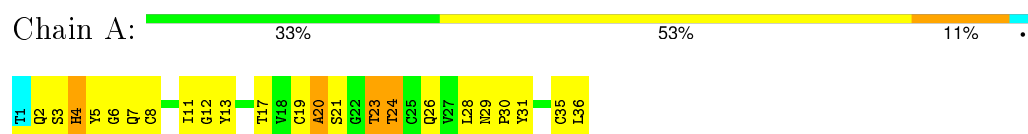
4.2.29 Score per residue for model 29

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.30 Score per residue for model 30

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.31 Score per residue for model 31

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



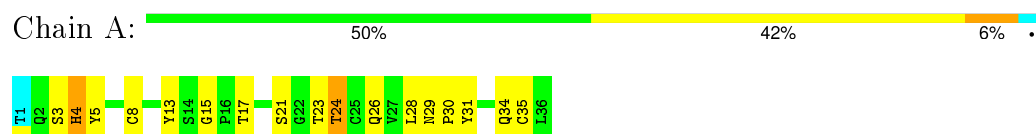
4.2.32 Score per residue for model 32 (medoid)

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



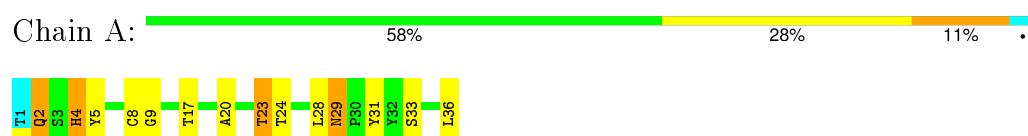
4.2.33 Score per residue for model 33

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



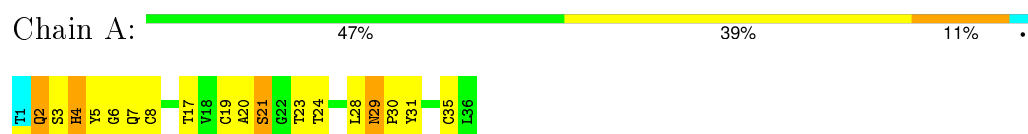
4.2.34 Score per residue for model 34

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



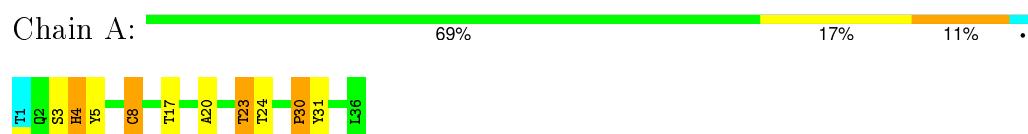
4.2.35 Score per residue for model 35

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.36 Score per residue for model 36

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.37 Score per residue for model 37

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.38 Score per residue for model 38

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



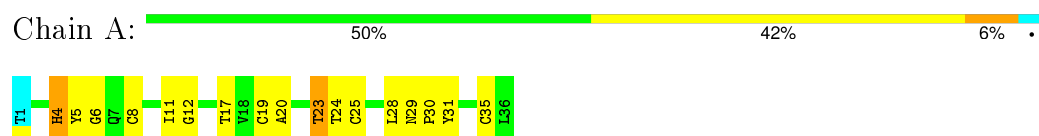
4.2.39 Score per residue for model 39

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



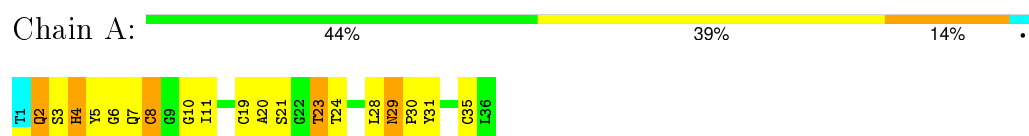
4.2.40 Score per residue for model 40

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



4.2.41 Score per residue for model 41

- Molecule 1: C-TERMINAL DOMAIN OF CELLOBIOHYDROLASE I



5 Refinement protocol and experimental data overview

Of the ? calculated structures, 41 were deposited, based on the following criterion: ?.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
X-PLOR	refinement	

No chemical shift data was provided. No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality [i](#)

6.1 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 (average) 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	1.06±0.01	1±0/259 (0.4±0.0%)	0.95±0.02	0±0/351 (0.0±0.0%)
All	All	1.06	41/10619 (0.4%)	0.95	0/14391 (0.0%)

All unique bond outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	4	HIS	CG-ND1	-6.03	1.25	1.38	21	41

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	253	226	226	10±3
All	All	10373	9266	9266	414

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 21.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:8:CYS:O	1:A:17:THR:HG22	0.80	1.76	38	26

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:26:GLN:HE22	1:A:36:LEU:H	0.74	1.25	15	2
1:A:2:GLN:O	1:A:20:ALA:HB2	0.68	1.88	24	6
1:A:4:HIS:CE1	1:A:5:TYR:CZ	0.66	2.83	27	27
1:A:4:HIS:CE1	1:A:5:TYR:CE2	0.65	2.84	2	30
1:A:4:HIS:ND1	1:A:5:TYR:CE2	0.64	2.66	21	21
1:A:4:HIS:NE2	1:A:5:TYR:CZ	0.63	2.66	5	4
1:A:2:GLN:HE22	1:A:7:GLN:N	0.63	1.91	3	1
1:A:26:GLN:NE2	1:A:36:LEU:H	0.62	1.93	15	2
1:A:23:THR:HG22	1:A:36:LEU:O	0.61	1.95	29	2
1:A:2:GLN:NE2	1:A:7:GLN:N	0.61	2.47	3	1
1:A:36:LEU:N	1:A:36:LEU:HD22	0.59	2.12	29	5
1:A:4:HIS:CG	1:A:5:TYR:CE2	0.59	2.90	5	9
1:A:11:ILE:HD12	1:A:30:PRO:O	0.59	1.97	9	20
1:A:8:CYS:SG	1:A:35:CYS:SG	0.58	3.01	7	18
1:A:26:GLN:NE2	1:A:36:LEU:N	0.58	2.51	10	2
1:A:2:GLN:HE22	1:A:8:CYS:N	0.58	1.97	30	2
1:A:23:THR:HG22	1:A:36:LEU:OXT	0.57	1.99	28	3
1:A:4:HIS:CD2	1:A:5:TYR:CE2	0.56	2.94	5	4
1:A:4:HIS:CD2	1:A:5:TYR:H	0.55	2.19	41	6
1:A:2:GLN:N	1:A:2:GLN:OE1	0.55	2.40	10	1
1:A:2:GLN:NE2	1:A:7:GLN:O	0.54	2.40	3	5
1:A:30:PRO:O	1:A:31:TYR:CD1	0.53	2.61	18	8
1:A:26:GLN:N	1:A:26:GLN:CD	0.53	2.62	20	3
1:A:2:GLN:O	1:A:20:ALA:N	0.52	2.42	34	4
1:A:2:GLN:HE22	1:A:7:GLN:C	0.52	2.08	16	3
1:A:2:GLN:NE2	1:A:13:TYR:OH	0.52	2.42	5	5
1:A:8:CYS:SG	1:A:19:CYS:SG	0.52	3.08	6	1
1:A:34:GLN:OE1	1:A:36:LEU:HD21	0.52	2.05	9	1
1:A:26:GLN:OE1	1:A:26:GLN:N	0.51	2.44	2	4
1:A:26:GLN:HE22	1:A:36:LEU:N	0.51	1.98	10	1
1:A:7:GLN:O	1:A:13:TYR:CZ	0.51	2.63	32	3
1:A:36:LEU:HD22	1:A:36:LEU:N	0.51	2.20	25	4
1:A:32:TYR:OH	1:A:34:GLN:NE2	0.51	2.43	11	1
1:A:32:TYR:OH	1:A:34:GLN:OE1	0.49	2.31	2	6
1:A:24:THR:C	1:A:26:GLN:HE22	0.49	2.11	20	2
1:A:26:GLN:CD	1:A:26:GLN:N	0.49	2.66	1	3
1:A:4:HIS:C	1:A:6:GLY:N	0.49	2.66	40	4
1:A:26:GLN:NE2	1:A:36:LEU:O	0.48	2.46	2	1
1:A:19:CYS:SG	1:A:25:CYS:N	0.48	2.87	21	2
1:A:30:PRO:O	1:A:31:TYR:CG	0.48	2.66	41	11
1:A:30:PRO:C	1:A:31:TYR:CD1	0.47	2.88	29	12

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:20:ALA:O	1:A:23:THR:OG1	0.46	2.34	28	20
1:A:9:GLY:O	1:A:33:SER:N	0.46	2.48	34	1
1:A:19:CYS:SG	1:A:24:THR:C	0.46	2.94	25	7
1:A:4:HIS:NE2	1:A:5:TYR:CE2	0.46	2.83	25	1
1:A:29:ASN:C	1:A:31:TYR:H	0.46	2.15	27	31
1:A:30:PRO:C	1:A:31:TYR:CG	0.45	2.89	22	5
1:A:2:GLN:HE22	1:A:7:GLN:CA	0.45	2.24	3	1
1:A:28:LEU:HD23	1:A:29:ASN:HD21	0.45	1.72	24	5
1:A:2:GLN:OE1	1:A:7:GLN:O	0.45	2.34	37	2
1:A:2:GLN:O	1:A:20:ALA:CB	0.45	2.64	31	1
1:A:13:TYR:CZ	1:A:15:GLY:CA	0.45	3.00	33	1
1:A:4:HIS:O	1:A:6:GLY:N	0.45	2.50	40	2
1:A:34:GLN:CG	1:A:35:CYS:N	0.45	2.80	10	1
1:A:24:THR:O	1:A:26:GLN:NE2	0.44	2.51	2	1
1:A:2:GLN:NE2	1:A:7:GLN:C	0.44	2.70	16	1
1:A:28:LEU:HD23	1:A:29:ASN:ND2	0.44	2.27	19	3
1:A:10:GLY:O	1:A:11:ILE:C	0.43	2.57	12	4
1:A:4:HIS:O	1:A:5:TYR:HB2	0.43	2.13	28	1
1:A:19:CYS:SG	1:A:25:CYS:CA	0.43	3.06	21	1
1:A:32:TYR:OH	1:A:34:GLN:CD	0.43	2.57	39	2
1:A:24:THR:O	1:A:26:GLN:OE1	0.43	2.36	1	5
1:A:20:ALA:O	1:A:21:SER:C	0.43	2.57	35	3
1:A:32:TYR:CD1	1:A:33:SER:N	0.43	2.87	25	1
1:A:8:CYS:O	1:A:17:THR:CG2	0.42	2.62	30	2
1:A:26:GLN:OE1	1:A:36:LEU:N	0.42	2.52	2	3
1:A:29:ASN:O	1:A:31:TYR:N	0.42	2.53	27	7
1:A:2:GLN:NE2	1:A:6:GLY:C	0.42	2.73	30	1
1:A:11:ILE:HG22	1:A:12:GLY:N	0.42	2.29	30	3
1:A:19:CYS:O	1:A:20:ALA:C	0.42	2.58	41	3
1:A:5:TYR:O	1:A:34:GLN:OE1	0.42	2.37	7	2
1:A:19:CYS:O	1:A:20:ALA:O	0.42	2.38	30	1
1:A:11:ILE:CD1	1:A:30:PRO:O	0.42	2.67	22	2
1:A:11:ILE:CG2	1:A:12:GLY:N	0.42	2.83	30	2
1:A:4:HIS:O	1:A:5:TYR:CB	0.41	2.68	28	1
1:A:2:GLN:OE1	1:A:7:GLN:C	0.41	2.59	41	1
1:A:26:GLN:OE1	1:A:34:GLN:O	0.41	2.39	33	1
1:A:13:TYR:CE2	1:A:15:GLY:O	0.41	2.73	33	1
1:A:3:SER:O	1:A:4:HIS:C	0.40	2.60	29	1
1:A:28:LEU:CG	1:A:29:ASN:ND2	0.40	2.84	3	1
1:A:14:SER:O	1:A:14:SER:OG	0.40	2.39	12	1
1:A:4:HIS:CD2	1:A:5:TYR:CD2	0.40	3.09	9	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:29:ASN:CB	1:A:30:PRO:CD	0.40	3.00	14	1
1:A:6:GLY:O	1:A:35:CYS:N	0.40	2.54	35	2
1:A:30:PRO:O	1:A:31:TYR:CD2	0.40	2.75	37	1

6.3 Torsion angles [i](#)

6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	34/36 (94%)	28±1 (82±4%)	6±1 (17±4%)	0±0 (1±1%)	26	73
All	All	1394/1476 (94%)	1140 (82%)	242 (17%)	12 (1%)	26	73

All 4 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	30	PRO	8
1	A	20	ALA	2
1	A	11	ILE	1
1	A	4	HIS	1

6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	28/29 (97%)	23±1 (83±5%)	5±1 (17±5%)	6	42
All	All	1148/1189 (97%)	952 (83%)	196 (17%)	6	42

All 11 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	24	THR	41
1	A	23	THR	40
1	A	3	SER	37
1	A	28	LEU	22
1	A	29	ASN	16
1	A	8	CYS	13
1	A	21	SER	11
1	A	2	GLN	9
1	A	16	PRO	4
1	A	34	GLN	2
1	A	26	GLN	1

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided