



Full wwPDB NMR Structure Validation Report ⓘ

Mar 8, 2026 – 07:18 AM UTC

PDB ID : 6EZ4 / pdb_00006ez4
BMRB ID : 34200
Title : NMR structure of the C-terminal domain of the human RPAP3 protein
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Deposited on : 2017-11-14

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

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

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
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

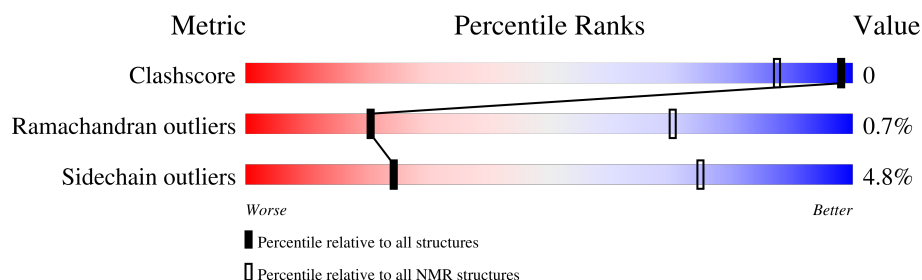
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 92%.

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	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

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	135	

2 Ensemble composition and analysis ⓘ

This entry contains 20 models. Model 6 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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:542-A:665 (124)	0.38	6

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 2 clusters and 3 single-model clusters were found.

Cluster number	Models
1	5, 6, 7, 9, 11, 12, 14, 15, 16, 17, 19
2	1, 2, 3, 4, 13, 20
Single-model clusters	8; 10; 18

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2250 atoms, of which 1137 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called RNA polymerase II-associated protein 3.

Mol	Chain	Residues	Atoms						Trace
1	A	135	Total	C	H	N	O	S	0
			2250	724	1137	182	202	5	

There are 4 discrepancies between the modelled and reference sequences:

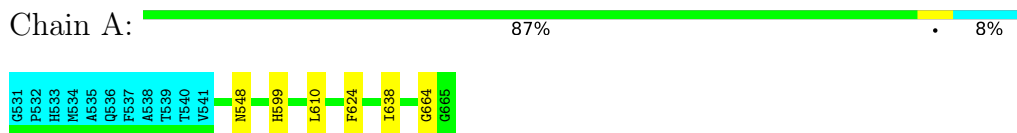
Chain	Residue	Modelled	Actual	Comment	Reference
A	531	GLY	-	expression tag	UNP Q9H6T3
A	532	PRO	-	expression tag	UNP Q9H6T3
A	533	HIS	-	expression tag	UNP Q9H6T3
A	534	MET	-	expression tag	UNP Q9H6T3

4 Residue-property plots [i](#)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: RNA polymerase II-associated protein 3

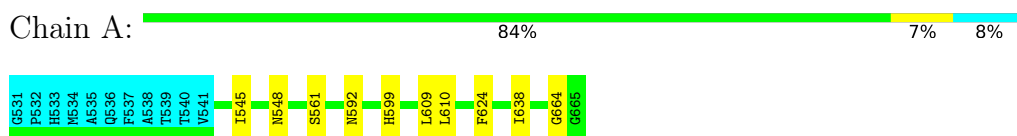


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

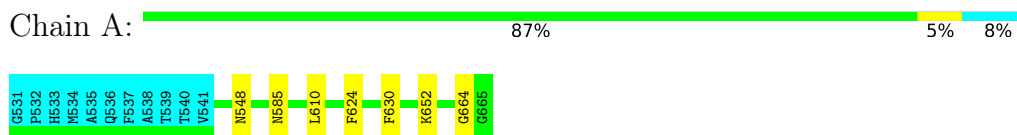
4.2.1 Score per residue for model 1

- Molecule 1: RNA polymerase II-associated protein 3



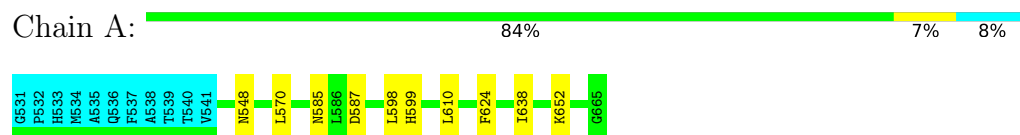
4.2.2 Score per residue for model 2

- Molecule 1: RNA polymerase II-associated protein 3



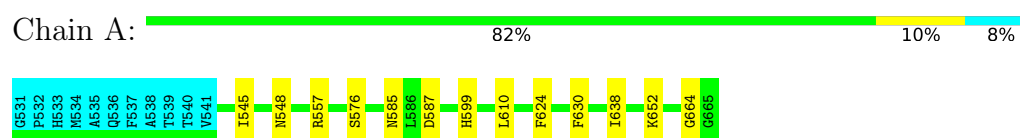
4.2.3 Score per residue for model 3

- Molecule 1: RNA polymerase II-associated protein 3



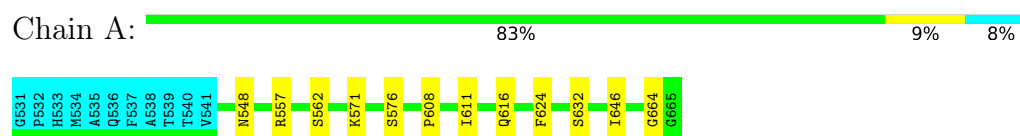
4.2.4 Score per residue for model 4

- Molecule 1: RNA polymerase II-associated protein 3



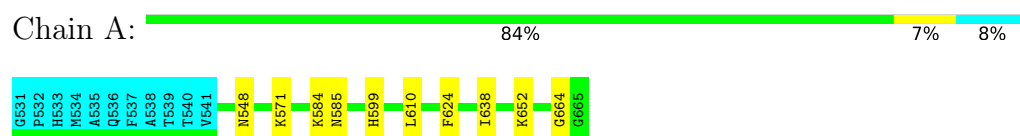
4.2.5 Score per residue for model 5

- Molecule 1: RNA polymerase II-associated protein 3



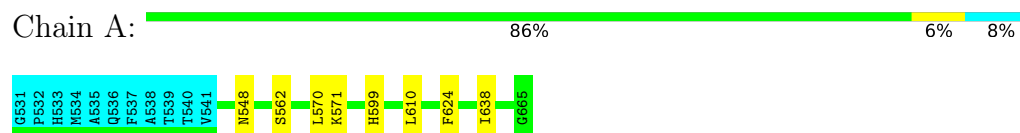
4.2.6 Score per residue for model 6 (medoid)

- Molecule 1: RNA polymerase II-associated protein 3



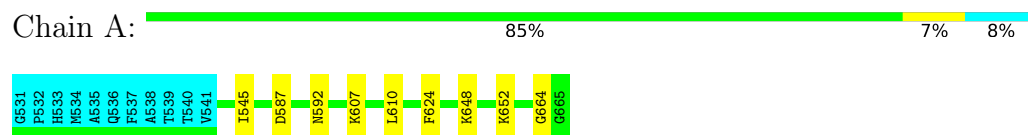
4.2.7 Score per residue for model 7

- Molecule 1: RNA polymerase II-associated protein 3



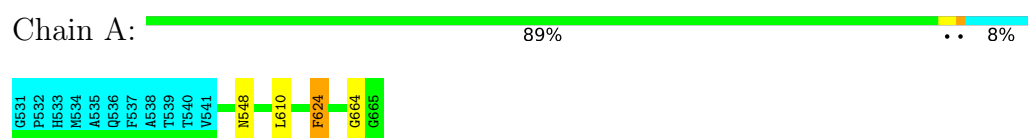
4.2.8 Score per residue for model 8

- Molecule 1: RNA polymerase II-associated protein 3



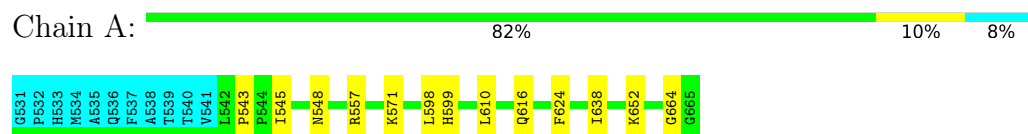
4.2.9 Score per residue for model 9

- Molecule 1: RNA polymerase II-associated protein 3



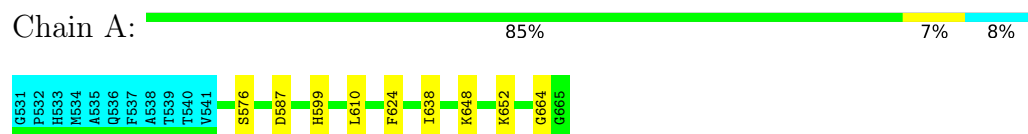
4.2.10 Score per residue for model 10

- Molecule 1: RNA polymerase II-associated protein 3



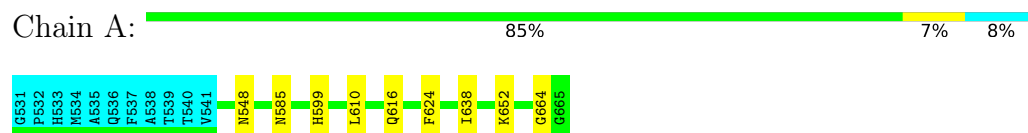
4.2.11 Score per residue for model 11

- Molecule 1: RNA polymerase II-associated protein 3



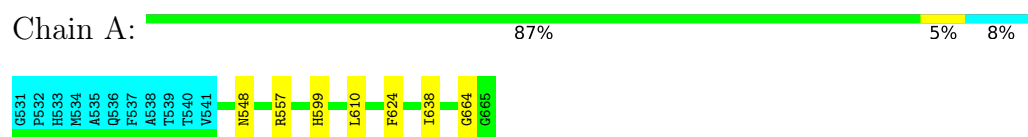
4.2.12 Score per residue for model 12

- Molecule 1: RNA polymerase II-associated protein 3



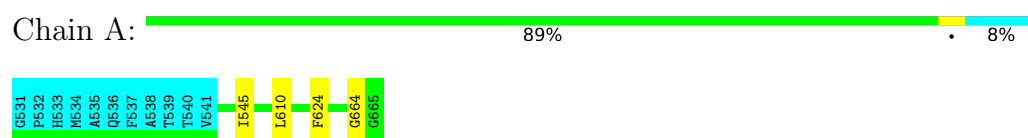
4.2.13 Score per residue for model 13

- Molecule 1: RNA polymerase II-associated protein 3



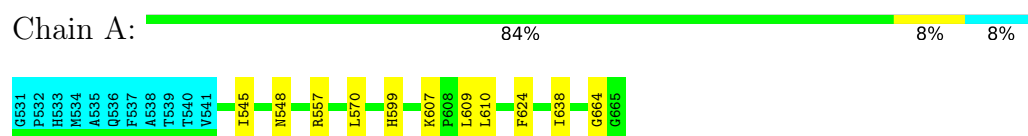
4.2.14 Score per residue for model 14

- Molecule 1: RNA polymerase II-associated protein 3



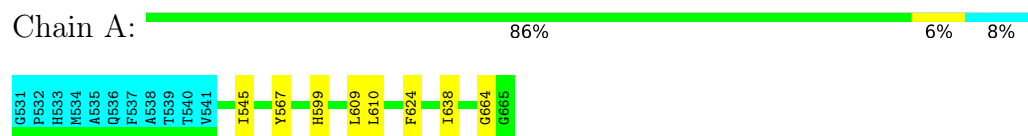
4.2.15 Score per residue for model 15

- Molecule 1: RNA polymerase II-associated protein 3



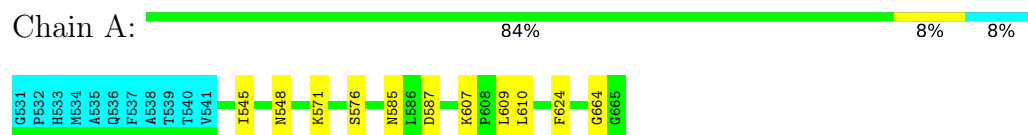
4.2.16 Score per residue for model 16

- Molecule 1: RNA polymerase II-associated protein 3



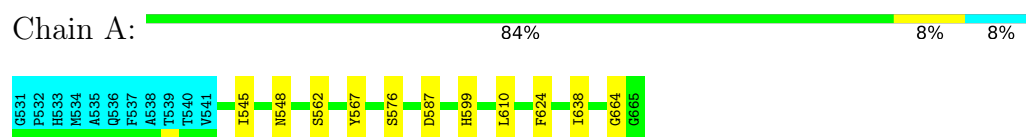
4.2.17 Score per residue for model 17

- Molecule 1: RNA polymerase II-associated protein 3



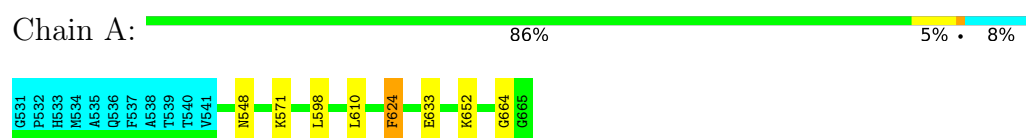
4.2.18 Score per residue for model 18

- Molecule 1: RNA polymerase II-associated protein 3



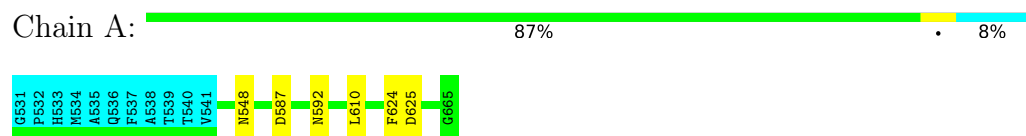
4.2.19 Score per residue for model 19

- Molecule 1: RNA polymerase II-associated protein 3



4.2.20 Score per residue for model 20

- Molecule 1: RNA polymerase II-associated protein 3



5 Refinement protocol and experimental data overview

The models were refined using the following method: *molecular dynamics*.

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest restraint energies*.

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

Software name	Classification	Version
CYANA	structure calculation	
TALOS	structure calculation	
Amber	refinement	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	1798
Number of shifts mapped to atoms	1798
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	92%

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	0.77±0.01	0±0/1056 (0.0± 0.0%)	1.31±0.02	2±2/1416 (0.1± 0.1%)
All	All	0.77	0/21120 (0.0%)	1.31	33/28320 (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	Planarity
1	A	0.0±0.0	0.3±0.5
All	All	0	7

There are no bond-length outliers.

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	607	LYS	CA-C-N	5.78	125.47	119.05	8	3
1	A	607	LYS	C-N-CA	5.78	125.47	119.05	8	3
1	A	598	LEU	CA-C-N	5.67	128.14	120.38	10	3
1	A	598	LEU	C-N-CA	5.67	128.14	120.38	10	3
1	A	548	ASN	CA-C-N	5.56	128.00	120.38	10	1
1	A	548	ASN	C-N-CA	5.56	128.00	120.38	10	1
1	A	543	PRO	N-CA-CB	5.52	106.09	103.22	10	1
1	A	545	ILE	CA-C-N	5.51	125.48	120.03	14	9
1	A	545	ILE	C-N-CA	5.51	125.48	120.03	14	9

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	A	585	ASN	Peptide	3
1	A	557	ARG	Sidechain	2
1	A	567	TYR	Sidechain	2

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	1033	1059	1059	1±1
All	All	20660	21180	21180	17

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 0.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:599:HIS:CG	1:A:638:ILE:HG23	0.49	2.43	4	10
1:A:599:HIS:CD2	1:A:638:ILE:HG23	0.44	2.48	12	3
1:A:624:PHE:C	1:A:624:PHE:CD1	0.44	2.95	19	1
1:A:624:PHE:CD1	1:A:624:PHE:C	0.42	2.97	9	1
1:A:608:PRO:HA	1:A:611:ILE:HD12	0.41	1.92	5	1
1:A:611:ILE:HG22	1:A:646:ILE:HD11	0.40	1.92	5	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	123/135 (91%)	120±1 (98±1%)	2±1 (2±1%)	1±0 (1±0%)	20	70
All	All	2460/2700 (91%)	2406 (98%)	37 (2%)	17 (1%)	20	70

All 1 unique Ramachandran outliers are listed below.

Mol	Chain	Res	Type	Models (Total)
1	A	664	GLY	17

6.3.2 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 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	117/125 (94%)	111±1 (95±1%)	6±1 (5±1%)	24	75
All	All	2340/2500 (94%)	2228 (95%)	112 (5%)	24	75

All 21 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	624	PHE	20
1	A	610	LEU	19
1	A	548	ASN	15
1	A	652	LYS	9
1	A	587	ASP	7
1	A	571	LYS	6
1	A	576	SER	5
1	A	609	LEU	4
1	A	592	ASN	3
1	A	585	ASN	3
1	A	570	LEU	3
1	A	557	ARG	3
1	A	562	SER	3
1	A	616	GLN	3
1	A	630	PHE	2
1	A	648	LYS	2
1	A	561	SER	1
1	A	632	SER	1
1	A	584	LYS	1
1	A	633	GLU	1
1	A	625	ASP	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 oligosaccharides 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

The completeness of assignment taking into account all chemical shift lists is 92% for the well-defined parts and 92% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: 535665_ref_dss_298K.str

7.1.1 Bookkeeping

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	1798
Number of shifts mapped to atoms	1798
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	134	-0.51 ± 0.16	Should be checked
$^{13}\text{C}_\beta$	131	0.25 ± 0.10	None needed (< 0.5 ppm)
$^{13}\text{C}'$	133	-0.27 ± 0.09	None needed (< 0.5 ppm)
^{15}N	125	0.53 ± 0.16	Should be applied

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 92%, i.e. 1679 atoms were assigned a chemical shift out of a possible 1818. 0 out of 21 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	606/607 (100%)	243/243 (100%)	247/248 (100%)	116/116 (100%)
Sidechain	980/1062 (92%)	665/687 (97%)	299/335 (89%)	16/40 (40%)

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	Total	¹H	¹³C	¹⁵N
Aromatic	93/149 (62%)	63/73 (86%)	30/74 (41%)	0/2 (0%)
Overall	1679/1818 (92%)	971/1003 (97%)	576/657 (88%)	132/158 (84%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 92%, i.e. 1793 atoms were assigned a chemical shift out of a possible 1954. 0 out of 22 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	654/661 (99%)	262/265 (99%)	267/270 (99%)	125/126 (99%)
Sidechain	1044/1127 (93%)	709/731 (97%)	318/355 (90%)	17/41 (41%)
Aromatic	95/166 (57%)	65/82 (79%)	30/81 (37%)	0/3 (0%)
Overall	1793/1954 (92%)	1036/1078 (96%)	615/706 (87%)	142/170 (84%)

7.1.4 Statistically unusual chemical shifts ⓘ

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	619	SER	HA	1.47	2.50 – 6.44	-7.6

7.1.5 Random Coil Index (RCI) plots ⓘ

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:

