



Full wwPDB NMR Structure Validation Report ⓘ

Mar 8, 2026 – 09:57 AM UTC

PDB ID : 4ASW / pdb_00004asw
BMRB ID : 18342
Title : Structure of the complex between the N-terminal dimerisation domain of Sgt2 and the UBL domain of Get5
Authors : Simon, A.C.; Simpson, P.J.; Goldstone, R.M.; Kryzstofinska, E.M.; Murray, J.W.; High, S.; Isaacson, R.L.
Deposited on : 2012-05-03

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

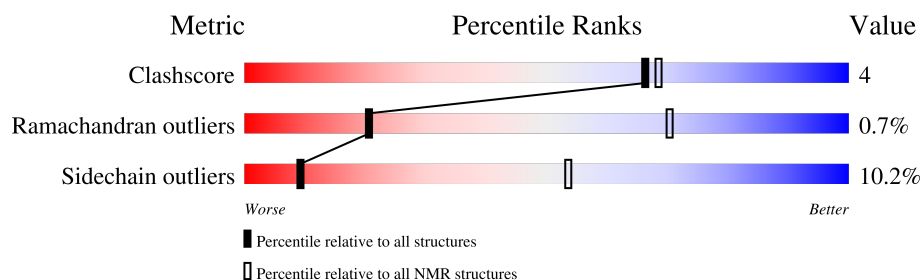
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 30%.

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	92	
1	B	92	
2	C	83	

2 Ensemble composition and analysis

This entry contains 9 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:16-A:92, B:14-B:90 (154)	0.63	6
2	C:19-C:95 (77)	0.48	7

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 3 clusters. No single-model clusters were found.

Cluster number	Models
1	1, 2, 5, 7, 8
2	3, 9
3	4, 6

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2223 atoms, of which 402 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2.

Mol	Chain	Residues	Atoms						Trace
1	A	79	Total	C	H	N	O	S	0
			723	369	129	96	125	4	
1	B	79	Total	C	H	N	O	S	0
			723	369	129	96	125	4	

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q12118
A	2	ALA	-	expression tag	UNP Q12118
A	3	HIS	-	expression tag	UNP Q12118
A	4	HIS	-	expression tag	UNP Q12118
A	5	HIS	-	expression tag	UNP Q12118
A	6	HIS	-	expression tag	UNP Q12118
A	7	HIS	-	expression tag	UNP Q12118
A	8	HIS	-	expression tag	UNP Q12118
A	9	VAL	-	expression tag	UNP Q12118
A	10	ASP	-	expression tag	UNP Q12118
A	11	ASP	-	expression tag	UNP Q12118
A	12	ASP	-	expression tag	UNP Q12118
A	13	ASP	-	expression tag	UNP Q12118
A	14	MET	-	expression tag	UNP Q12118
B	1	MET	-	expression tag	UNP Q12118
B	2	ALA	-	expression tag	UNP Q12118
B	3	HIS	-	expression tag	UNP Q12118
B	4	HIS	-	expression tag	UNP Q12118
B	5	HIS	-	expression tag	UNP Q12118
B	6	HIS	-	expression tag	UNP Q12118
B	7	HIS	-	expression tag	UNP Q12118
B	8	HIS	-	expression tag	UNP Q12118
B	9	VAL	-	expression tag	UNP Q12118
B	10	ASP	-	expression tag	UNP Q12118
B	11	ASP	-	expression tag	UNP Q12118
B	12	ASP	-	expression tag	UNP Q12118
B	13	ASP	-	expression tag	UNP Q12118
B	14	MET	-	expression tag	UNP Q12118

- Molecule 2 is a protein called UBIQUITIN-LIKE PROTEIN MDY2.

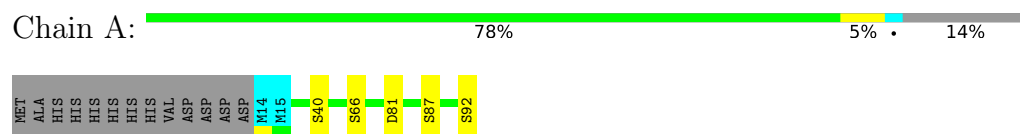
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
2	C	81	777	409	144	107	116	1	0

4 Residue-property plots [i](#)

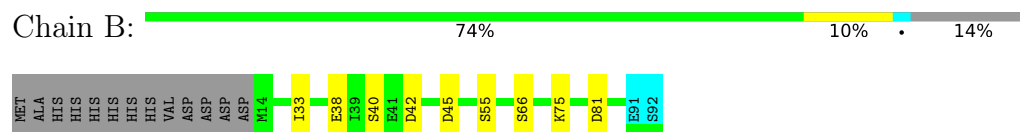
4.1 Average score per residue in the NMR ensemble

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.

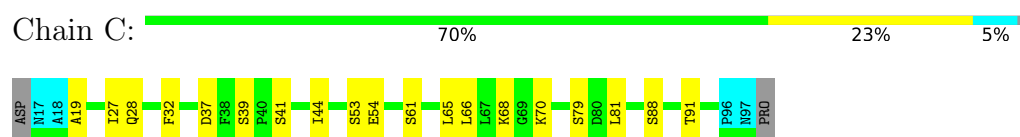
- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2

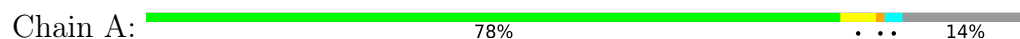


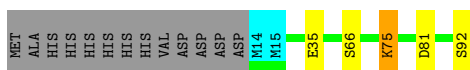
4.2 Scores per residue for each member of the ensemble

Colouring as in section [4.1](#) above.

4.2.1 Score per residue for model 1

- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2





- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



4.2.2 Score per residue for model 2

- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



4.2.3 Score per residue for model 3

- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2

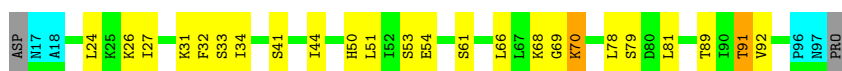




- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



4.2.4 Score per residue for model 4

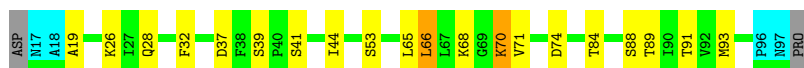
- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



4.2.5 Score per residue for model 5

- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2





- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



4.2.6 Score per residue for model 6 (medoid)

- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



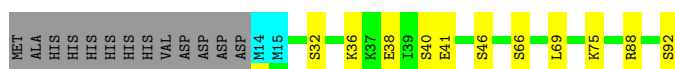
- Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



4.2.7 Score per residue for model 7

- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2





- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



4.2.8 Score per residue for model 8

- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



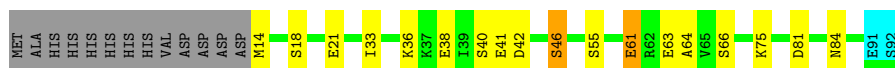
4.2.9 Score per residue for model 9

- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2





- Molecule 1: SMALL GLUTAMINE-RICH TETRATRICOPEPTIDE REPEAT-CONTAINING PROTEIN 2



- Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



5 Refinement protocol and experimental data overview

The models were refined using the following method: *ARIA*.

Of the 100 calculated structures, 9 were deposited, based on the following criterion: *LOWEST ENERGY*.

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

Software name	Classification	Version
ARIA	refinement	
NMRView	structure solution	

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	938
Number of shifts mapped to atoms	497
Number of unparsed shifts	0
Number of shifts with mapping errors	441
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	30%

6 Model quality [i](#)

6.1 Standard geometry [i](#)

There are no covalent bond-length or bond-angle outliers.

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
2	C	1.0±0.0	0.0±0.0
All	All	9	0

There are no bond-length outliers.

There are no bond-angle outliers.

All unique chiral outliers are listed below.

Mol	Chain	Res	Type	Atoms	Models (Total)
2	C	19	ALA	CA	9

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	578	125	563	4±2
1	B	578	126	569	4±1
2	C	608	137	652	8±2
All	All	15876	3492	16056	125

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:49:VAL:HG21	2:C:69:GLY:HA2	0.69	1.63	3	1
1:A:49:VAL:HG11	2:C:66:LEU:HD21	0.68	1.66	4	1
2:C:66:LEU:HG	2:C:91:THR:HB	0.67	1.65	9	1
1:A:90:PRO:HB3	1:B:75:LYS:HD3	0.67	1.67	4	1
2:C:70:LYS:HA	2:C:70:LYS:HE3	0.66	1.66	6	3
1:B:45:ASP:HB3	2:C:27:ILE:HB	0.65	1.67	8	1
1:B:33:ILE:HG23	1:B:38:GLU:HB2	0.64	1.68	7	4
1:B:33:ILE:HG23	1:B:38:GLU:HB3	0.64	1.67	4	4
1:A:87:SER:HA	1:B:14:MET:HB2	0.61	1.70	4	1
2:C:19:ALA:HB1	2:C:37:ASP:HB3	0.59	1.72	9	4
2:C:45:LEU:HG	2:C:74:ASP:HB3	0.59	1.75	2	1
2:C:27:ILE:H	2:C:27:ILE:HD13	0.58	1.58	1	1
1:A:87:SER:HG	1:B:14:MET:N	0.57	1.98	9	1
1:B:21:GLU:HG2	1:B:72:SER:HB3	0.57	1.76	5	1
2:C:64:LYS:HB3	2:C:93:MET:HB2	0.57	1.76	6	1
1:B:78:HIS:HB2	1:B:81:ASP:OD2	0.56	2.00	5	1
1:A:75:LYS:HD3	1:B:90:PRO:HB3	0.55	1.78	5	1
1:A:83:LEU:HD21	1:B:25:LEU:HD12	0.55	1.78	8	1
2:C:23:THR:HB	2:C:89:THR:HG23	0.55	1.78	7	1
2:C:66:LEU:HB2	2:C:91:THR:HB	0.54	1.79	5	2
1:B:18:SER:HB2	1:B:21:GLU:HG3	0.54	1.79	1	2
2:C:26:LYS:HD3	2:C:32:PHE:CE2	0.53	2.38	4	1
2:C:44:ILE:HG23	2:C:65:LEU:HD12	0.53	1.79	5	5
1:A:49:VAL:HG21	2:C:66:LEU:HD13	0.53	1.81	9	1
2:C:58:SER:HB2	2:C:62:GLU:OE1	0.52	2.04	8	1
2:C:78:LEU:HA	2:C:81:LEU:HG	0.52	1.81	6	4
1:A:45:ASP:OD2	2:C:70:LYS:HA	0.52	2.05	3	1
2:C:26:LYS:HE3	2:C:32:PHE:CE2	0.52	2.40	5	1
1:A:32:SER:O	1:A:36:LYS:HG2	0.51	2.05	5	4
2:C:25:LYS:HA	2:C:32:PHE:O	0.51	2.06	2	2
1:B:51:MET:SD	1:B:62:ARG:HD3	0.51	2.45	6	1
2:C:28:GLN:NE2	2:C:93:MET:HA	0.51	2.21	4	1
2:C:44:ILE:HB	2:C:74:ASP:HA	0.51	1.82	4	1
1:A:45:ASP:HB3	2:C:71:VAL:H	0.50	1.67	9	1
2:C:84:THR:HB	2:C:85:PRO:HD2	0.50	1.84	6	2
2:C:23:THR:HG23	2:C:35:GLU:HG2	0.49	1.83	6	3
2:C:28:GLN:HE22	2:C:93:MET:HA	0.49	1.66	4	1
2:C:50:HIS:O	2:C:54:GLU:HG2	0.49	2.07	6	3
2:C:44:ILE:HD11	2:C:81:LEU:HD21	0.49	1.84	2	2
1:A:33:ILE:HG23	1:A:38:GLU:HB2	0.49	1.85	2	1
1:B:42:ASP:HA	2:C:31:LYS:HD3	0.48	1.84	1	1
2:C:24:LEU:HD13	2:C:51:LEU:HD21	0.47	1.84	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:61:GLU:HB3	1:B:64:ALA:HB3	0.47	1.87	9	1
2:C:26:LYS:HG2	2:C:28:GLN:O	0.46	2.10	5	2
1:A:69:LEU:O	1:A:75:LYS:HB3	0.46	2.11	7	1
1:B:55:SER:HA	1:B:60:PHE:CE2	0.46	2.46	2	4
2:C:59:HIS:HB2	2:C:62:GLU:HG3	0.45	1.86	6	1
2:C:66:LEU:HG	2:C:71:VAL:HA	0.45	1.87	4	1
2:C:29:ALA:HA	2:C:30:PRO:C	0.44	2.37	2	2
1:A:75:LYS:HD2	1:A:75:LYS:H	0.44	1.71	1	1
1:B:21:GLU:HB3	1:B:74:PHE:HE2	0.44	1.73	2	1
1:A:55:SER:HA	1:A:60:PHE:CE2	0.44	2.48	5	3
2:C:26:LYS:HA	2:C:92:VAL:O	0.44	2.12	3	1
1:A:84:ASN:O	1:B:14:MET:HA	0.44	2.12	4	1
2:C:59:HIS:HB2	2:C:62:GLU:OE2	0.44	2.12	8	1
2:C:84:THR:HB	2:C:85:PRO:CD	0.44	2.43	6	2
2:C:27:ILE:O	2:C:31:LYS:HE2	0.44	2.13	9	2
2:C:19:ALA:HB2	2:C:37:ASP:HB3	0.43	1.89	4	1
1:B:46:SER:HA	2:C:27:ILE:HD12	0.43	1.89	9	1
1:B:14:MET:HB3	1:B:73:GLU:OE1	0.43	2.14	6	1
1:A:49:VAL:CG2	2:C:69:GLY:HA2	0.43	2.41	3	1
2:C:32:PHE:HE1	2:C:34:ILE:HG12	0.43	1.74	3	1
2:C:45:LEU:O	2:C:49:GLN:HG3	0.43	2.13	8	1
2:C:26:LYS:HE2	2:C:94:ILE:HD12	0.43	1.90	5	1
2:C:62:GLU:O	2:C:94:ILE:HA	0.42	2.14	1	1
2:C:48:LYS:HD3	2:C:60:ILE:HG23	0.42	1.88	5	1
2:C:54:GLU:O	2:C:56:LYS:HE2	0.42	2.13	8	1
1:B:49:VAL:HG21	2:C:91:THR:HG21	0.42	1.90	1	1
2:C:66:LEU:HD23	2:C:71:VAL:HA	0.42	1.91	5	1
1:B:79:LEU:O	1:B:83:LEU:HB2	0.42	2.15	8	1
1:A:88:ARG:HD3	1:B:77:GLN:OE1	0.41	2.15	5	1
2:C:62:GLU:HB3	2:C:95:LYS:O	0.41	2.14	8	1
1:B:63:GLU:OE1	1:B:63:GLU:HA	0.41	2.14	2	1
1:A:17:ALA:HB2	1:B:84:ASN:HB2	0.41	1.92	9	1
1:A:29:TYR:O	1:A:33:ILE:HG13	0.41	2.14	3	2
1:A:41:GLU:H	1:A:41:GLU:CD	0.41	2.23	7	1
1:A:79:LEU:O	1:A:83:LEU:HB2	0.41	2.16	8	1
1:A:25:LEU:HD12	1:B:83:LEU:HD21	0.41	1.92	3	1
2:C:39:SER:O	2:C:42:ASP:HB2	0.41	2.16	8	1
1:A:22:ILE:HG12	1:B:83:LEU:HD23	0.41	1.92	3	1
1:A:90:PRO:HG3	1:B:75:LYS:HD3	0.41	1.92	8	1
1:A:49:VAL:HB	2:C:66:LEU:HD13	0.41	1.92	3	1
2:C:24:LEU:O	2:C:33:SER:HA	0.41	2.16	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:53:CYS:HB3	1:B:46:SER:HB3	0.40	1.92	2	1
1:A:87:SER:HA	1:B:14:MET:SD	0.40	2.56	3	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	76/92 (83%)	73±1 (96±1%)	3±1 (4±1%)	0±0 (0±0%)	49	83
1	B	76/92 (83%)	72±1 (94±2%)	4±1 (5±2%)	1±1 (1±1%)	13	61
2	C	77/83 (93%)	74±1 (96±2%)	3±1 (4±2%)	1±1 (1±1%)	20	70
All	All	2061/2403 (86%)	1964 (95%)	83 (4%)	14 (1%)	20	70

All 6 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
2	C	19	ALA	4
1	B	16	SER	4
1	B	17	ALA	3
1	B	15	MET	1
1	A	16	SER	1
2	C	69	GLY	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	63/77 (82%)	58±2 (91±3%)	5±2 (9±3%)	12	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	63/77 (82%)	56±2 (90±4%)	7±2 (10±4%)	9	53
2	C	72/77 (94%)	64±2 (89±3%)	8±2 (11±3%)	8	51
All	All	1782/2079 (86%)	1601 (90%)	181 (10%)	9	53

All 67 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	66	SER	8
2	C	41	SER	8
1	B	66	SER	7
2	C	39	SER	7
2	C	70	LYS	7
2	C	79	SER	7
2	C	88	SER	7
1	A	81	ASP	5
1	A	92	SER	5
1	B	42	ASP	5
2	C	61	SER	5
2	C	68	LYS	5
1	A	40	SER	5
1	B	40	SER	5
2	C	53	SER	5
1	B	45	ASP	4
1	B	81	ASP	4
1	B	85	SER	4
1	B	75	LYS	4
1	B	52	ASP	3
1	B	61	GLU	3
1	A	38	GLU	3
1	B	73	GLU	3
1	A	18	SER	3
2	C	66	LEU	3
1	A	35	GLU	2
2	C	28	GLN	2
1	A	87	SER	2
2	C	45	LEU	2
1	A	42	ASP	2
1	B	63	GLU	2
2	C	89	THR	2
2	C	91	THR	2

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Mol	Chain	Res	Type	Models (Total)
1	B	18	SER	2
1	A	63	GLU	2
1	B	46	SER	2
1	A	46	SER	2
1	A	88	ARG	2
1	B	71	LYS	2
1	A	75	LYS	1
2	C	27	ILE	1
2	C	75	ASN	1
2	C	80	ASP	1
1	A	52	ASP	1
1	B	37	LYS	1
1	B	62	ARG	1
2	C	74	ASP	1
1	A	37	LYS	1
1	A	55	SER	1
2	C	84	THR	1
2	C	43	THR	1
2	C	54	GLU	1
1	A	16	SER	1
1	A	48	ASN	1
1	B	35	GLU	1
1	B	48	ASN	1
2	C	87	ASN	1
1	B	56	GLU	1
2	C	23	THR	1
2	C	95	LYS	1
1	A	41	GLU	1
1	A	71	LYS	1
1	B	83	LEU	1
1	B	36	LYS	1
1	B	41	GLU	1
1	B	55	SER	1
2	C	83	VAL	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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 30% for the well-defined parts and 30% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: 4ASW

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	938
Number of shifts mapped to atoms	497
Number of unparsed shifts	0
Number of shifts with mapping errors	441
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- No matching atom found in the structure. All 441 occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	15	MET	HA	4.469	0.03	1
1	A	15	MET	HB2	2.01	0.03	2
1	A	15	MET	HB3	2.041	0.03	2
1	A	15	MET	HG2	2.476	0.03	2
1	A	15	MET	HG3	2.569	0.03	2
1	A	16	SER	HA	4.439	0.03	1
1	A	16	SER	HB2	3.796	0.03	1
1	A	16	SER	HB3	3.796	0.03	1
1	A	17	ALA	HA	4.454	0.03	1
1	A	17	ALA	HB1	1.291	0.03	1
1	A	17	ALA	HB2	1.291	0.03	1
1	A	17	ALA	HB3	1.291	0.03	1
1	A	18	SER	HA	4.607	0.03	1
1	A	18	SER	HB2	4.019	0.03	2

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	18	SER	HB3	4.43	0.03	2
1	A	19	LYS	HA	4.086	0.03	1
1	A	19	LYS	HB2	1.91	0.03	2
1	A	19	LYS	HB3	2.096	0.03	2
1	A	19	LYS	HD2	1.98	0.03	2
1	A	19	LYS	HD3	2.0	0.03	2
1	A	19	LYS	HE2	3.172	0.03	2
1	A	19	LYS	HE3	3.232	0.03	2
1	A	19	LYS	HG2	1.688	0.03	2
1	A	19	LYS	HG3	1.804	0.03	2
1	A	20	GLU	HA	3.831	0.03	1
1	A	20	GLU	HB2	2.279	0.03	2
1	A	20	GLU	HB3	1.986	0.03	2
1	A	20	GLU	HG2	2.456	0.03	2
1	A	20	GLU	HG3	2.644	0.03	2
1	A	21	GLU	HA	3.637	0.03	1
1	A	21	GLU	HB2	2.521	0.03	2
1	A	21	GLU	HB3	1.846	0.03	2
1	A	21	GLU	HG2	2.081	0.03	2
1	A	21	GLU	HG3	2.355	0.03	2
1	A	22	ILE	HA	3.196	0.03	1
1	A	22	ILE	HB	1.712	0.03	1
1	A	22	ILE	HD11	0.66	0.03	1
1	A	22	ILE	HD12	0.66	0.03	1
1	A	22	ILE	HD13	0.66	0.03	1
1	A	22	ILE	HG12	0.535	0.03	1
1	A	22	ILE	HG13	1.861	0.03	1
1	A	22	ILE	HG21	0.162	0.03	1
1	A	22	ILE	HG22	0.162	0.03	1
1	A	22	ILE	HG23	0.162	0.03	1
1	A	23	ALA	HA	3.851	0.03	1
1	A	23	ALA	HB1	1.031	0.03	1
1	A	23	ALA	HB2	1.031	0.03	1
1	A	23	ALA	HB3	1.031	0.03	1
1	A	24	ALA	HA	3.874	0.03	1
1	A	24	ALA	HB1	1.149	0.03	1
1	A	24	ALA	HB2	1.149	0.03	1
1	A	24	ALA	HB3	1.149	0.03	1
1	A	25	LEU	HA	3.827	0.03	1
1	A	25	LEU	HB2	1.376	0.03	2
1	A	25	LEU	HB3	1.673	0.03	2

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	25	LEU	HD11	-0.07	0.03	2
1	A	25	LEU	HD12	-0.07	0.03	2
1	A	25	LEU	HD13	-0.07	0.03	2
1	A	25	LEU	HD21	0.428	0.03	2
1	A	25	LEU	HD22	0.428	0.03	2
1	A	25	LEU	HD23	0.428	0.03	2
1	A	25	LEU	HG	1.227	0.03	1
1	A	26	ILE	HA	3.237	0.03	1
1	A	26	ILE	HB	2.159	0.03	1
1	A	26	ILE	HD11	0.451	0.03	1
1	A	26	ILE	HD12	0.451	0.03	1
1	A	26	ILE	HD13	0.451	0.03	1
1	A	26	ILE	HG12	0.3	0.03	1
1	A	26	ILE	HG13	1.565	0.03	1
1	A	26	ILE	HG21	0.757	0.03	1
1	A	26	ILE	HG22	0.757	0.03	1
1	A	26	ILE	HG23	0.757	0.03	1
1	A	27	VAL	HA	3.615	0.03	1
1	A	27	VAL	HB	2.441	0.03	1
1	A	27	VAL	HG11	1.368	0.03	2
1	A	27	VAL	HG12	1.368	0.03	2
1	A	27	VAL	HG13	1.368	0.03	2
1	A	27	VAL	HG21	1.172	0.03	2
1	A	27	VAL	HG22	1.172	0.03	2
1	A	27	VAL	HG23	1.172	0.03	2
1	A	28	ASN	HA	4.524	0.03	1
1	A	28	ASN	HB2	2.71	0.03	2
1	A	28	ASN	HB3	2.75	0.03	2
1	A	29	TYR	HA	4.314	0.03	1
1	A	29	TYR	HB2	3.086	0.03	2
1	A	29	TYR	HB3	2.97	0.03	2
1	A	29	TYR	HD1	6.665	0.03	3
1	A	29	TYR	HD2	6.665	0.03	3
1	A	29	TYR	HE1	6.62	0.03	3
1	A	29	TYR	HE2	6.62	0.03	3
1	A	30	PHE	HA	4.439	0.03	1
1	A	30	PHE	HB2	3.167	0.03	2
1	A	30	PHE	HB3	3.552	0.03	2
1	A	30	PHE	HD1	7.189	0.03	3
1	A	30	PHE	HD2	7.189	0.03	3
1	A	30	PHE	HE1	7.316	0.03	3

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	30	PHE	HE2	7.316	0.03	3
1	A	31	SER	HA	4.175	0.03	1
1	A	31	SER	HB2	4.071	0.03	1
1	A	31	SER	HB3	4.071	0.03	1
1	A	32	SER	HA	4.289	0.03	1
1	A	32	SER	HB2	3.993	0.03	2
1	A	32	SER	HB3	4.064	0.03	2
1	A	33	ILE	HA	3.767	0.03	1
1	A	33	ILE	HB	2.081	0.03	1
1	A	33	ILE	HD11	0.728	0.03	1
1	A	33	ILE	HD12	0.728	0.03	1
1	A	33	ILE	HD13	0.728	0.03	1
1	A	33	ILE	HG12	1.145	0.03	1
1	A	33	ILE	HG13	1.36	0.03	1
1	A	33	ILE	HG21	0.728	0.03	1
1	A	33	ILE	HG22	0.728	0.03	1
1	A	33	ILE	HG23	0.728	0.03	1
1	A	34	VAL	HA	3.569	0.03	1
1	A	34	VAL	HB	2.212	0.03	1
1	A	34	VAL	HG11	0.995	0.03	2
1	A	34	VAL	HG12	0.995	0.03	2
1	A	34	VAL	HG13	0.995	0.03	2
1	A	34	VAL	HG21	0.956	0.03	2
1	A	34	VAL	HG22	0.956	0.03	2
1	A	34	VAL	HG23	0.956	0.03	2
1	A	35	GLU	HA	3.999	0.03	1
1	A	35	GLU	HB2	2.151	0.03	2
1	A	35	GLU	HB3	2.088	0.03	2
1	A	35	GLU	HG2	2.273	0.03	2
1	A	35	GLU	HG3	2.37	0.03	2
1	A	36	LYS	HA	4.184	0.03	1
1	A	36	LYS	HB2	1.728	0.03	2
1	A	36	LYS	HB3	2.037	0.03	2
1	A	36	LYS	HD2	1.62	0.03	2
1	A	36	LYS	HD3	1.59	0.03	2
1	A	36	LYS	HE2	2.913	0.03	2
1	A	36	LYS	HE3	2.911	0.03	2
1	A	36	LYS	HG2	1.423	0.03	2
1	A	36	LYS	HG3	1.642	0.03	2
1	A	37	LYS	HA	3.93	0.03	1
1	A	37	LYS	HB2	1.987	0.03	2

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	37	LYS	HB3	2.081	0.03	2
1	A	37	LYS	HD2	1.712	0.03	2
1	A	37	LYS	HD3	1.752	0.03	2
1	A	37	LYS	HE2	3.036	0.03	2
1	A	37	LYS	HE3	3.028	0.03	2
1	A	37	LYS	HG2	1.399	0.03	2
1	A	37	LYS	HG3	1.38	0.03	2
1	A	38	GLU	HA	4.305	0.03	1
1	A	38	GLU	HB2	1.846	0.03	2
1	A	38	GLU	HB3	2.245	0.03	2
1	A	38	GLU	HG2	2.2	0.03	2
1	A	38	GLU	HG3	2.186	0.03	2
1	A	39	ILE	HA	4.446	0.03	1
1	A	39	ILE	HB	1.55	0.03	1
1	A	39	ILE	HD11	0.219	0.03	1
1	A	39	ILE	HD12	0.219	0.03	1
1	A	39	ILE	HD13	0.219	0.03	1
1	A	39	ILE	HG12	0.984	0.03	1
1	A	39	ILE	HG13	1.431	0.03	1
1	A	39	ILE	HG21	0.585	0.03	1
1	A	39	ILE	HG22	0.585	0.03	1
1	A	39	ILE	HG23	0.585	0.03	1
1	A	40	SER	HA	4.325	0.03	1
1	A	40	SER	HB2	4.282	0.03	2
1	A	40	SER	HB3	4.039	0.03	2
1	A	41	GLU	HA	3.976	0.03	1
1	A	41	GLU	HB2	2.057	0.03	2
1	A	41	GLU	HB3	2.049	0.03	2
1	A	41	GLU	HG2	2.363	0.03	1
1	A	41	GLU	HG3	2.363	0.03	1
1	A	42	ASP	HA	4.409	0.03	1
1	A	42	ASP	HB2	2.48	0.03	2
1	A	42	ASP	HB3	2.645	0.03	2
1	A	43	GLY	HA2	3.503	0.03	2
1	A	43	GLY	HA3	3.637	0.03	2
1	A	44	ALA	HA	3.71	0.03	1
1	A	44	ALA	HB1	1.447	0.03	1
1	A	44	ALA	HB2	1.447	0.03	1
1	A	44	ALA	HB3	1.447	0.03	1
1	A	45	ASP	HA	4.446	0.03	1
1	A	45	ASP	HB2	2.744	0.03	2

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	45	ASP	HB3	2.911	0.03	2
1	A	46	SER	HA	4.203	0.03	1
1	A	46	SER	HB2	3.882	0.03	2
1	A	46	SER	HB3	3.96	0.03	2
1	A	47	LEU	HA	3.818	0.03	1
1	A	47	LEU	HB2	1.051	0.03	2
1	A	47	LEU	HB3	1.62	0.03	2
1	A	47	LEU	HD11	0.42	0.03	2
1	A	47	LEU	HD12	0.42	0.03	2
1	A	47	LEU	HD13	0.42	0.03	2
1	A	47	LEU	HD21	0.135	0.03	2
1	A	47	LEU	HD22	0.135	0.03	2
1	A	47	LEU	HD23	0.135	0.03	2
1	A	47	LEU	HG	1.443	0.03	1
1	A	48	ASN	HA	4.478	0.03	1
1	A	48	ASN	HB2	2.977	0.03	2
1	A	48	ASN	HB3	3.039	0.03	2
1	A	49	VAL	HA	3.843	0.03	1
1	A	49	VAL	HB	2.41	0.03	1
1	A	49	VAL	HG11	1.062	0.03	2
1	A	49	VAL	HG12	1.062	0.03	2
1	A	49	VAL	HG13	1.062	0.03	2
1	A	49	VAL	HG21	0.91	0.03	2
1	A	49	VAL	HG22	0.91	0.03	2
1	A	49	VAL	HG23	0.91	0.03	2
1	A	50	ALA	HA	3.827	0.03	1
1	A	50	ALA	HB1	1.486	0.03	1
1	A	50	ALA	HB2	1.486	0.03	1
1	A	50	ALA	HB3	1.486	0.03	1
1	A	51	MET	HA	3.97	0.03	1
1	A	51	MET	HB2	1.924	0.03	2
1	A	51	MET	HB3	2.644	0.03	2
1	A	51	MET	HE1	2.05	0.03	1
1	A	51	MET	HE2	2.05	0.03	1
1	A	51	MET	HE3	2.05	0.03	1
1	A	51	MET	HG2	2.558	0.03	2
1	A	51	MET	HG3	3.135	0.03	2
1	A	52	ASP	HA	4.531	0.03	1
1	A	52	ASP	HB2	2.759	0.03	2
1	A	52	ASP	HB3	3.044	0.03	2
1	A	53	CYS	HA	4.108	0.03	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	53	CYS	HB2	2.941	0.03	2
1	A	53	CYS	HB3	3.05	0.03	2
1	A	54	ILE	HA	3.263	0.03	1
1	A	54	ILE	HB	1.744	0.03	1
1	A	54	ILE	HD11	1.127	0.03	1
1	A	54	ILE	HD12	1.127	0.03	1
1	A	54	ILE	HD13	1.127	0.03	1
1	A	54	ILE	HG12	0.6	0.03	1
1	A	54	ILE	HG13	1.88	0.03	1
1	A	54	ILE	HG21	1.068	0.03	1
1	A	54	ILE	HG22	1.068	0.03	1
1	A	54	ILE	HG23	1.068	0.03	1
1	A	55	SER	HA	3.888	0.03	1
1	A	55	SER	HB2	2.958	0.03	2
1	A	55	SER	HB3	3.744	0.03	2
1	A	56	GLU	HA	3.909	0.03	1
1	A	56	GLU	HB2	1.971	0.03	2
1	A	56	GLU	HB3	2.065	0.03	2
1	A	56	GLU	HG2	2.269	0.03	1
1	A	56	GLU	HG3	2.269	0.03	1
1	A	57	ALA	HA	3.808	0.03	1
1	A	57	ALA	HB1	1.027	0.03	1
1	A	57	ALA	HB2	1.027	0.03	1
1	A	57	ALA	HB3	1.027	0.03	1
1	A	58	PHE	HA	4.472	0.03	1
1	A	58	PHE	HB2	2.669	0.03	2
1	A	58	PHE	HB3	3.48	0.03	2
1	A	58	PHE	HD1	7.64	0.03	3
1	A	58	PHE	HD2	7.64	0.03	3
1	A	58	PHE	HE1	7.16	0.03	3
1	A	58	PHE	HE2	7.16	0.03	3
1	A	59	GLY	HA2	3.874	0.03	2
1	A	59	GLY	HA3	4.039	0.03	2
1	A	60	PHE	HA	5.2	0.03	1
1	A	60	PHE	HB2	2.981	0.03	2
1	A	60	PHE	HB3	3.13	0.03	2
1	A	60	PHE	HD1	7.079	0.03	3
1	A	60	PHE	HD2	7.079	0.03	3
1	A	60	PHE	HE1	7.12	0.03	3
1	A	60	PHE	HE2	7.12	0.03	3
1	A	60	PHE	HZ	6.72	0.03	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	61	GLU	HA	4.591	0.03	1
1	A	61	GLU	HB2	1.959	0.03	2
1	A	61	GLU	HB3	2.359	0.03	2
1	A	61	GLU	HG2	2.35	0.03	2
1	A	61	GLU	HG3	2.38	0.03	2
1	A	62	ARG	HA	4.09	0.03	1
1	A	62	ARG	HB2	2.01	0.03	2
1	A	62	ARG	HB3	2.006	0.03	2
1	A	62	ARG	HD2	3.21	0.03	2
1	A	62	ARG	HD3	3.23	0.03	2
1	A	62	ARG	HG2	1.689	0.03	2
1	A	62	ARG	HG3	1.869	0.03	2
1	A	63	GLU	HA	4.238	0.03	1
1	A	63	GLU	HB2	2.087	0.03	2
1	A	63	GLU	HB3	2.13	0.03	2
1	A	63	GLU	HG2	2.33	0.03	2
1	A	63	GLU	HG3	2.4	0.03	2
1	A	64	ALA	HA	4.532	0.03	1
1	A	64	ALA	HB1	1.679	0.03	1
1	A	64	ALA	HB2	1.679	0.03	1
1	A	64	ALA	HB3	1.679	0.03	1
1	A	65	VAL	HA	3.258	0.03	1
1	A	65	VAL	HB	2.191	0.03	1
1	A	65	VAL	HG11	1.238	0.03	2
1	A	65	VAL	HG12	1.238	0.03	2
1	A	65	VAL	HG13	1.238	0.03	2
1	A	65	VAL	HG21	0.95	0.03	2
1	A	65	VAL	HG22	0.95	0.03	2
1	A	65	VAL	HG23	0.95	0.03	2
1	A	66	SER	HA	3.941	0.03	1
1	A	66	SER	HB2	3.919	0.03	1
1	A	66	SER	HB3	3.919	0.03	1
1	A	67	GLY	HA2	3.896	0.03	1
1	A	67	GLY	HA3	3.896	0.03	1
1	A	68	ILE	HA	3.557	0.03	1
1	A	68	ILE	HB	1.532	0.03	1
1	A	68	ILE	HD11	0.209	0.03	1
1	A	68	ILE	HD12	0.209	0.03	1
1	A	68	ILE	HD13	0.209	0.03	1
1	A	68	ILE	HG12	1.53	0.03	1
1	A	68	ILE	HG13	0.835	0.03	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	68	ILE	HG21	0.593	0.03	1
1	A	68	ILE	HG22	0.593	0.03	1
1	A	68	ILE	HG23	0.593	0.03	1
1	A	69	LEU	HA	4.067	0.03	1
1	A	69	LEU	HB2	1.398	0.03	2
1	A	69	LEU	HB3	1.774	0.03	2
1	A	69	LEU	HD11	0.747	0.03	2
1	A	69	LEU	HD12	0.747	0.03	2
1	A	69	LEU	HD13	0.747	0.03	2
1	A	69	LEU	HD21	0.812	0.03	2
1	A	69	LEU	HD22	0.812	0.03	2
1	A	69	LEU	HD23	0.812	0.03	2
1	A	69	LEU	HG	1.638	0.03	1
1	A	70	GLY	HA2	3.851	0.03	2
1	A	70	GLY	HA3	3.995	0.03	2
1	A	71	LYS	HA	4.56	0.03	1
1	A	71	LYS	HB2	1.818	0.03	2
1	A	71	LYS	HB3	1.979	0.03	2
1	A	71	LYS	HD2	1.65	0.03	1
1	A	71	LYS	HD3	1.65	0.03	1
1	A	71	LYS	HE2	2.956	0.03	1
1	A	71	LYS	HE3	2.956	0.03	1
1	A	71	LYS	HG2	1.532	0.03	2
1	A	71	LYS	HG3	1.489	0.03	2
1	A	72	SER	HA	4.767	0.03	1
1	A	72	SER	HB2	3.915	0.03	2
1	A	72	SER	HB3	4.101	0.03	2
1	A	73	GLU	HA	4.148	0.03	1
1	A	73	GLU	HB3	1.62	0.03	1
1	A	73	GLU	HG2	1.765	0.03	2
1	A	73	GLU	HG3	1.846	0.03	2
1	A	74	PHE	HA	4.613	0.03	1
1	A	74	PHE	HB2	2.63	0.03	2
1	A	74	PHE	HB3	3.311	0.03	2
1	A	74	PHE	HD1	6.93	0.03	3
1	A	74	PHE	HD2	6.93	0.03	3
1	A	74	PHE	HE1	7.01	0.03	3
1	A	74	PHE	HE2	7.01	0.03	3
1	A	74	PHE	HZ	6.87	0.03	1
1	A	75	LYS	HA	3.866	0.03	1
1	A	75	LYS	HB2	1.844	0.03	2

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	75	LYS	HB3	2.017	0.03	2
1	A	75	LYS	HD2	1.679	0.03	2
1	A	75	LYS	HD3	1.792	0.03	2
1	A	75	LYS	HE2	2.981	0.03	1
1	A	75	LYS	HE3	2.981	0.03	1
1	A	75	LYS	HG2	1.305	0.03	2
1	A	75	LYS	HG3	1.458	0.03	2
1	A	76	GLY	HA2	3.725	0.03	2
1	A	76	GLY	HA3	4.017	0.03	2
1	A	77	GLN	HA	4.582	0.03	1
1	A	77	GLN	HB2	1.97	0.03	2
1	A	77	GLN	HB3	1.99	0.03	2
1	A	77	GLN	HG2	2.1	0.03	2
1	A	77	GLN	HG3	2.237	0.03	2
1	A	78	HIS	HA	4.754	0.03	1
1	A	78	HIS	HB2	3.09	0.03	2
1	A	78	HIS	HB3	3.37	0.03	2
1	A	78	HIS	HD2	7.209	0.03	1
1	A	78	HIS	HE1	8.13	0.03	1
1	A	79	LEU	HA	3.58	0.03	1
1	A	79	LEU	HB2	1.112	0.03	1
1	A	79	LEU	HB3	1.308	0.03	1
1	A	79	LEU	HD11	0.381	0.03	2
1	A	79	LEU	HD12	0.381	0.03	2
1	A	79	LEU	HD13	0.381	0.03	2
1	A	79	LEU	HD21	-0.034	0.03	2
1	A	79	LEU	HD22	-0.034	0.03	2
1	A	79	LEU	HD23	-0.034	0.03	2
1	A	79	LEU	HG	0.898	0.03	1
1	A	80	ALA	HA	3.694	0.03	1
1	A	80	ALA	HB1	1.472	0.03	1
1	A	80	ALA	HB2	1.472	0.03	1
1	A	80	ALA	HB3	1.472	0.03	1
1	A	81	ASP	HA	4.447	0.03	1
1	A	81	ASP	HB2	2.77	0.03	2
1	A	81	ASP	HB3	2.942	0.03	2
1	A	82	ILE	HA	3.634	0.03	1
1	A	82	ILE	HB	1.556	0.03	1
1	A	82	ILE	HD11	0.538	0.03	1
1	A	82	ILE	HD12	0.538	0.03	1
1	A	82	ILE	HD13	0.538	0.03	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	82	ILE	HG12	0.937	0.03	1
1	A	82	ILE	HG13	1.642	0.03	1
1	A	82	ILE	HG21	0.687	0.03	1
1	A	82	ILE	HG22	0.687	0.03	1
1	A	82	ILE	HG23	0.687	0.03	1
1	A	83	LEU	HA	3.849	0.03	1
1	A	83	LEU	HB2	1.384	0.03	2
1	A	83	LEU	HB3	1.618	0.03	2
1	A	83	LEU	HD11	0.295	0.03	2
1	A	83	LEU	HD12	0.295	0.03	2
1	A	83	LEU	HD13	0.295	0.03	2
1	A	83	LEU	HD21	0.296	0.03	2
1	A	83	LEU	HD22	0.296	0.03	2
1	A	83	LEU	HD23	0.296	0.03	2
1	A	83	LEU	HG	1.513	0.03	1
1	A	84	ASN	HA	4.686	0.03	1
1	A	84	ASN	HB2	2.92	0.03	2
1	A	84	ASN	HB3	2.9	0.03	2
1	A	85	SER	HA	4.368	0.03	1
1	A	85	SER	HB2	3.992	0.03	2
1	A	85	SER	HB3	3.988	0.03	2
1	A	86	ALA	HA	4.289	0.03	1
1	A	86	ALA	HB1	1.429	0.03	1
1	A	86	ALA	HB2	1.429	0.03	1
1	A	86	ALA	HB3	1.429	0.03	1
1	A	87	SER	HA	4.382	0.03	1
1	A	87	SER	HB2	3.843	0.03	2
1	A	87	SER	HB3	3.839	0.03	2
1	A	88	ARG	HA	4.363	0.03	1
1	A	88	ARG	HB2	1.723	0.03	2
1	A	88	ARG	HB3	1.823	0.03	2
1	A	88	ARG	HD2	3.157	0.03	1
1	A	88	ARG	HD3	3.157	0.03	1
1	A	88	ARG	HG2	1.58	0.03	2
1	A	88	ARG	HG3	1.584	0.03	2
1	A	89	VAL	HA	4.348	0.03	1
1	A	89	VAL	HB	2.056	0.03	1
1	A	89	VAL	HG11	0.917	0.03	1
1	A	89	VAL	HG12	0.917	0.03	1
1	A	89	VAL	HG13	0.917	0.03	1
1	A	89	VAL	HG21	0.917	0.03	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	89	VAL	HG22	0.917	0.03	1
1	A	89	VAL	HG23	0.917	0.03	1
1	A	90	PRO	HA	4.407	0.03	1
1	A	90	PRO	HB2	1.932	0.03	2
1	A	90	PRO	HB3	2.28	0.03	2
1	A	90	PRO	HD2	3.842	0.03	2
1	A	90	PRO	HD3	3.638	0.03	2
1	A	90	PRO	HG2	1.944	0.03	2
1	A	90	PRO	HG3	2.022	0.03	2
1	A	91	GLU	HA	4.257	0.03	1
1	A	91	GLU	HB2	1.939	0.03	2
1	A	91	GLU	HB3	2.057	0.03	2
1	A	91	GLU	HG2	2.276	0.03	1
1	A	91	GLU	HG3	2.276	0.03	1
1	A	92	SER	HA	4.38	0.03	1
1	A	92	SER	HB2	3.82	0.03	1
1	A	92	SER	HB3	3.82	0.03	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
$^{13}\text{C}_\alpha$	78	-0.49 ± 0.14	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	73	0.26 ± 0.07	None needed (< 0.5 ppm)
$^{13}\text{C}'$	73	-0.28 ± 0.08	None needed (< 0.5 ppm)
^{15}N	76	0.51 ± 0.21	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 30%, i.e. 927 atoms were assigned a chemical shift out of a possible 3071. 0 out of 35 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	381/1156 (33%)	157/468 (34%)	149/462 (32%)	75/226 (33%)
Sidechain	498/1731 (29%)	340/1133 (30%)	152/553 (27%)	6/45 (13%)
Aromatic	48/184 (26%)	24/91 (26%)	24/79 (30%)	0/14 (0%)
Overall	927/3071 (30%)	521/1692 (31%)	325/1094 (30%)	81/285 (28%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 30%, i.e. 938 atoms were assigned a chemical shift out of a possible 3166. 0 out of 35 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	386/1194 (32%)	159/483 (33%)	151/478 (32%)	76/233 (33%)
Sidechain	504/1788 (28%)	344/1170 (29%)	154/571 (27%)	6/47 (13%)
Aromatic	48/184 (26%)	24/91 (26%)	24/79 (30%)	0/14 (0%)
Overall	938/3166 (30%)	527/1744 (30%)	329/1128 (29%)	82/294 (28%)

7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots [i](#)

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:

