



Full wwPDB NMR Structure Validation Report ⓘ

Mar 8, 2026 – 03:56 PM UTC

PDB ID : 7QUU / pdb_00007quu
BMRB ID : 34702
Title : Red1-Iss10 complex
Authors : Mackereth, C.D.; Kadlec, J.; Laroussi, H.
Deposited on : 2022-01-18

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

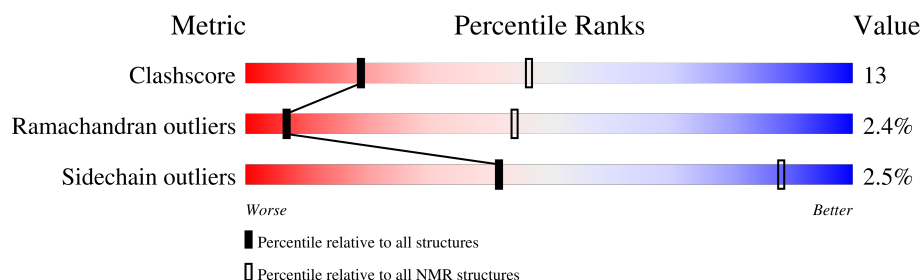
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 88%.

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	55	
2	B	49	

2 Ensemble composition and analysis

This entry contains 20 models. Model 15 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:4-A:43, B:193-B:234 (82)	0.61	15

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 2 single-model clusters were found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1497 atoms, of which 751 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Uncharacterized protein C7D4.14c.

Mol	Chain	Residues	Atoms					Trace
1	A	44	Total	C	H	N	O	0
			751	238	378	67	68	

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	initiating methionine	UNP O14269
A	-9	GLY	-	expression tag	UNP O14269
A	-8	TRP	-	expression tag	UNP O14269
A	-7	SER	-	expression tag	UNP O14269
A	-6	HIS	-	expression tag	UNP O14269
A	-5	PRO	-	expression tag	UNP O14269
A	-4	GLN	-	expression tag	UNP O14269
A	-3	PHE	-	expression tag	UNP O14269
A	-2	GLU	-	expression tag	UNP O14269
A	-1	LYS	-	expression tag	UNP O14269
A	0	SER	-	expression tag	UNP O14269
A	1	SER	-	expression tag	UNP O14269

- Molecule 2 is a protein called NURS complex subunit red1.

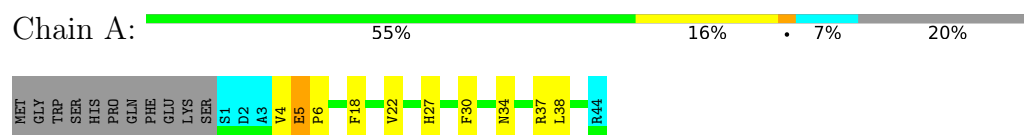
Mol	Chain	Residues	Atoms						Trace
2	B	49	Total	C	H	N	O	S	0
			746	237	373	61	74	1	

There are 5 discrepancies between the modelled and reference sequences:

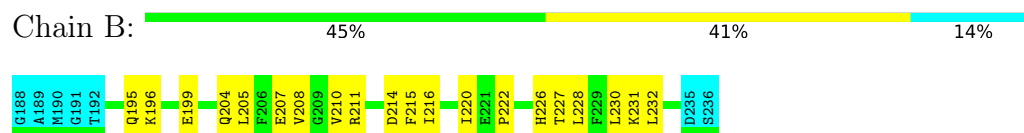
Chain	Residue	Modelled	Actual	Comment	Reference
B	188	GLY	-	expression tag	UNP Q9UTR8
B	189	ALA	-	expression tag	UNP Q9UTR8
B	190	MET	-	expression tag	UNP Q9UTR8
B	191	GLY	-	expression tag	UNP Q9UTR8
B	236	SER	-	expression tag	UNP Q9UTR8

4.2.2 Score per residue for model 2

- Molecule 1: Uncharacterized protein C7D4.14c

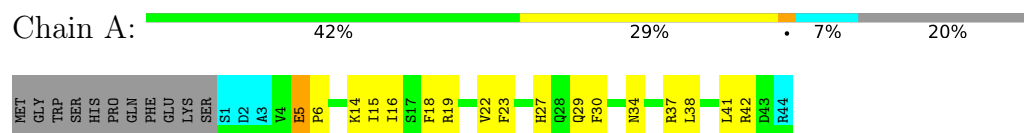


- Molecule 2: NURS complex subunit red1

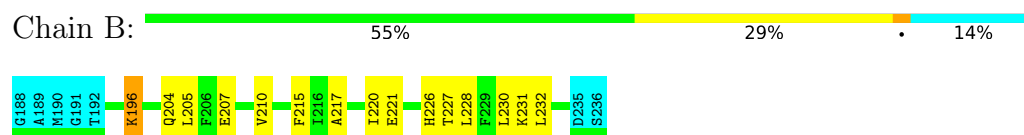


4.2.3 Score per residue for model 3

- Molecule 1: Uncharacterized protein C7D4.14c

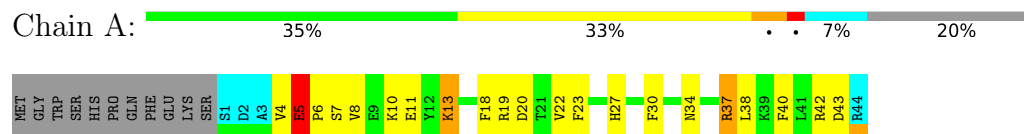


- Molecule 2: NURS complex subunit red1

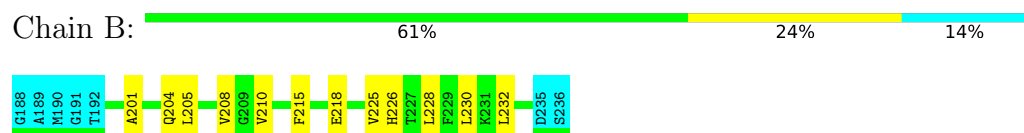


4.2.4 Score per residue for model 4

- Molecule 1: Uncharacterized protein C7D4.14c

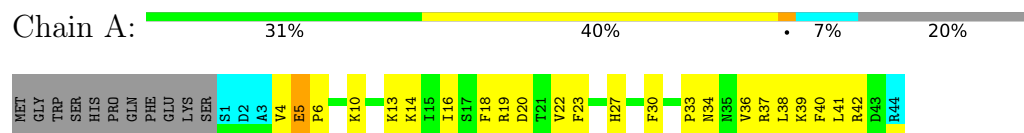


- Molecule 2: NURS complex subunit red1

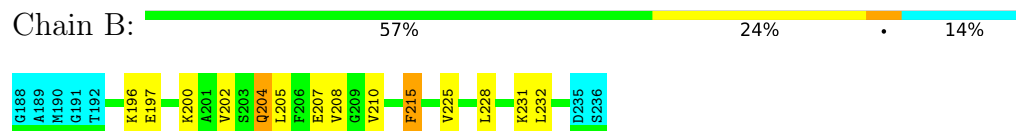


4.2.5 Score per residue for model 5

- Molecule 1: Uncharacterized protein C7D4.14c

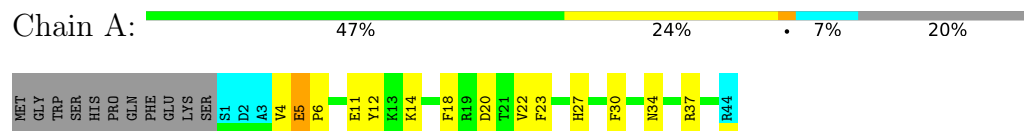


- Molecule 2: NURS complex subunit red1



4.2.6 Score per residue for model 6

- Molecule 1: Uncharacterized protein C7D4.14c

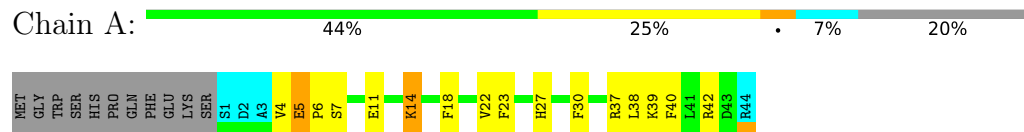


- Molecule 2: NURS complex subunit red1

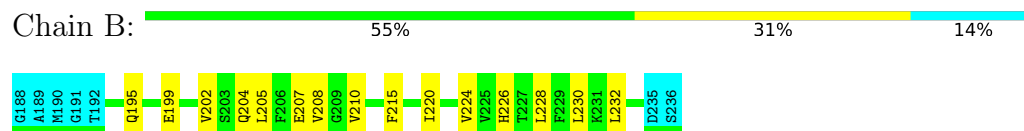


4.2.7 Score per residue for model 7

- Molecule 1: Uncharacterized protein C7D4.14c

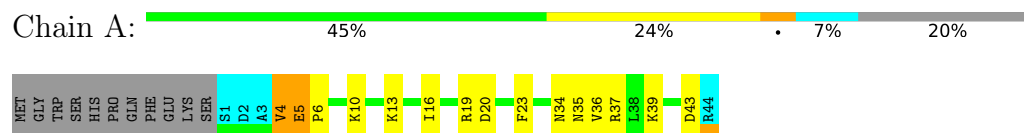


- Molecule 2: NURS complex subunit red1

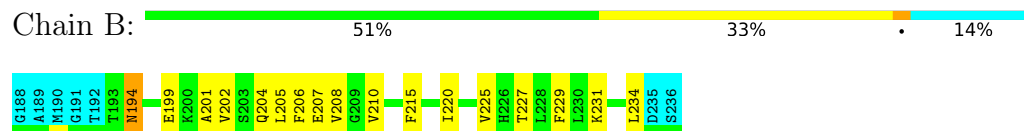


4.2.8 Score per residue for model 8

- Molecule 1: Uncharacterized protein C7D4.14c

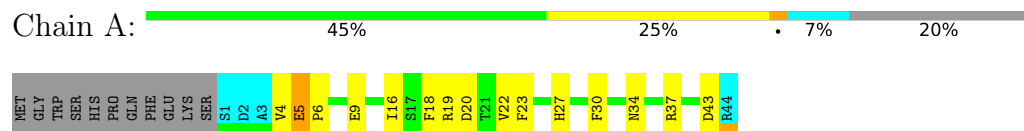


- Molecule 2: NURS complex subunit red1

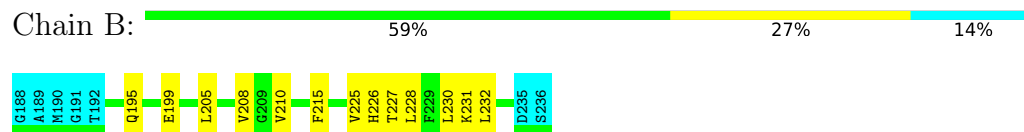


4.2.9 Score per residue for model 9

- Molecule 1: Uncharacterized protein C7D4.14c

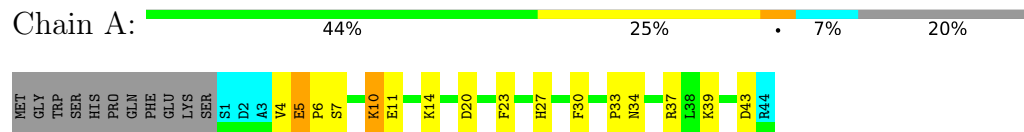


- Molecule 2: NURS complex subunit red1

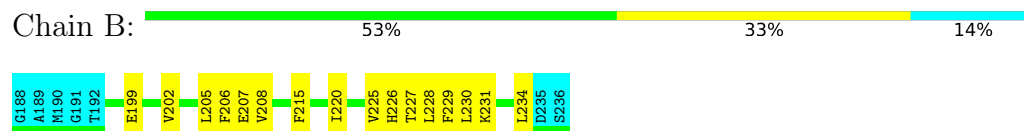


4.2.10 Score per residue for model 10

- Molecule 1: Uncharacterized protein C7D4.14c

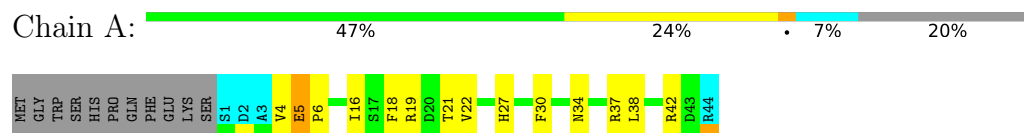


- Molecule 2: NURS complex subunit red1

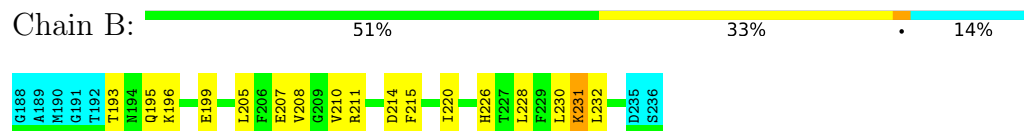


4.2.11 Score per residue for model 11

- Molecule 1: Uncharacterized protein C7D4.14c

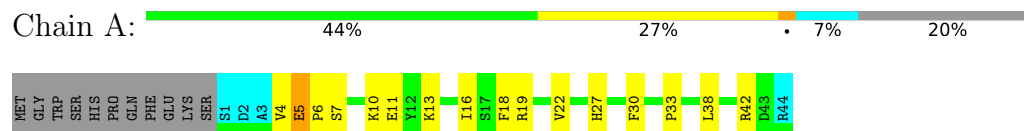


- Molecule 2: NURS complex subunit red1



4.2.12 Score per residue for model 12

- Molecule 1: Uncharacterized protein C7D4.14c

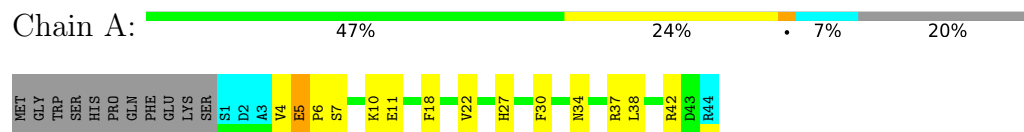


- Molecule 2: NURS complex subunit red1

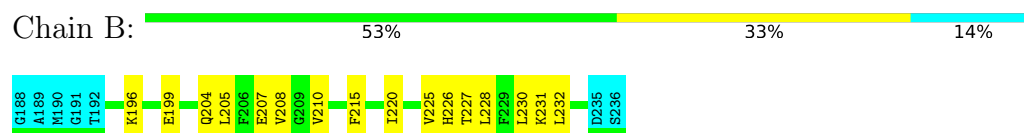


4.2.13 Score per residue for model 13

- Molecule 1: Uncharacterized protein C7D4.14c

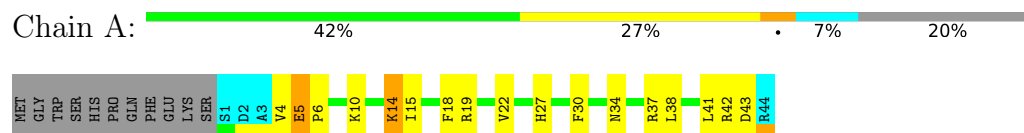


- Molecule 2: NURS complex subunit red1

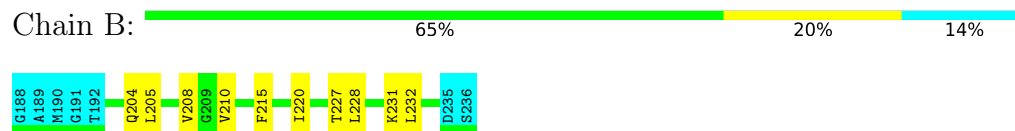


4.2.14 Score per residue for model 14

- Molecule 1: Uncharacterized protein C7D4.14c

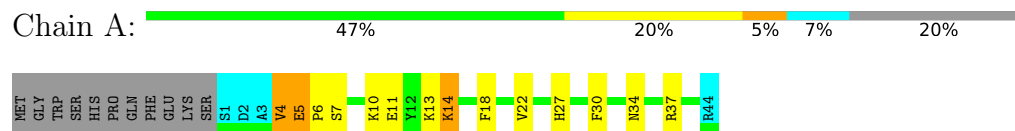


- Molecule 2: NURS complex subunit red1

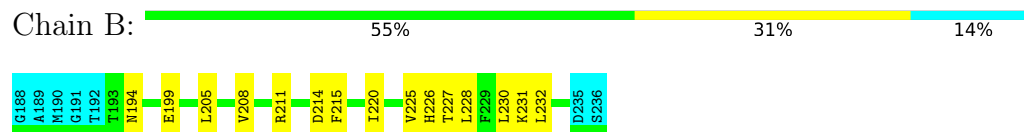


4.2.15 Score per residue for model 15 (medoid)

- Molecule 1: Uncharacterized protein C7D4.14c

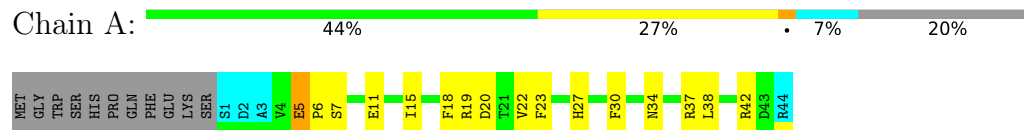


- Molecule 2: NURS complex subunit red1

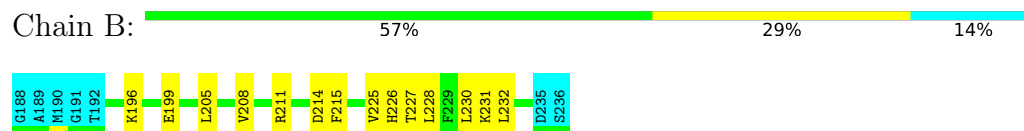


4.2.16 Score per residue for model 16

- Molecule 1: Uncharacterized protein C7D4.14c

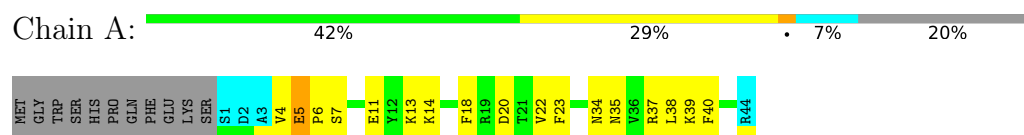


- Molecule 2: NURS complex subunit red1



4.2.17 Score per residue for model 17

- Molecule 1: Uncharacterized protein C7D4.14c

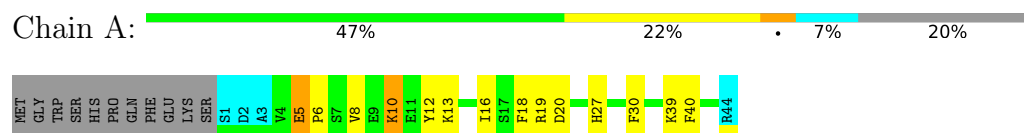


- Molecule 2: NURS complex subunit red1

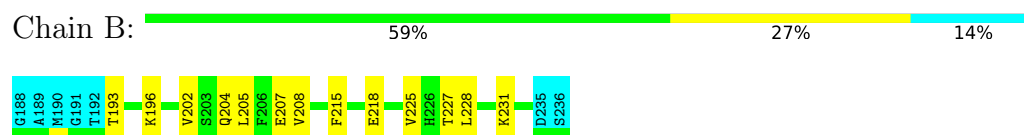


4.2.18 Score per residue for model 18

- Molecule 1: Uncharacterized protein C7D4.14c

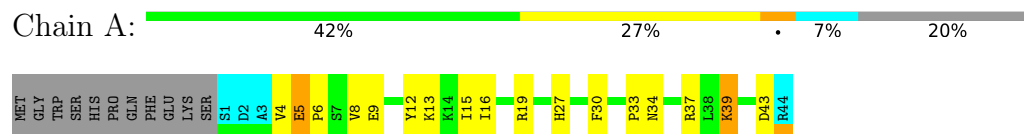


- Molecule 2: NURS complex subunit red1

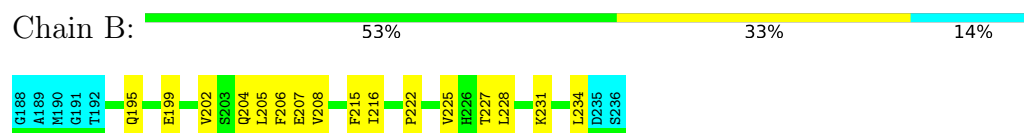


4.2.19 Score per residue for model 19

- Molecule 1: Uncharacterized protein C7D4.14c

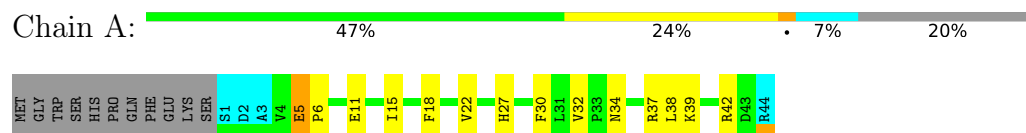


- Molecule 2: NURS complex subunit red1

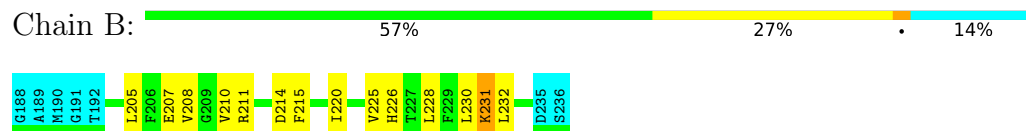


4.2.20 Score per residue for model 20

- Molecule 1: Uncharacterized protein C7D4.14c



- Molecule 2: NURS complex subunit red1



5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing*.

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

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

Software name	Classification	Version
ARIA	structure calculation	2.3

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	1117
Number of shifts mapped to atoms	1117
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	88%

6 Model quality [i](#)

6.1 Standard geometry [i](#)

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

There are no bond-length outliers.

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	342	351	350	9±3
2	B	330	335	334	9±2
All	All	13440	13720	13680	346

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 13.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:228:LEU:O	2:B:232:LEU:HG	0.72	1.84	3	15
1:A:5:GLU:HB2	1:A:6:PRO:CD	0.69	2.18	19	3
1:A:15:ILE:O	1:A:19:ARG:HG3	0.67	1.88	3	3
1:A:12:TYR:OH	2:B:194:ASN:HA	0.65	1.92	6	1
2:B:205:LEU:O	2:B:208:VAL:HG22	0.64	1.93	19	18
2:B:206:PHE:CD2	2:B:234:LEU:HD13	0.64	2.28	19	2
2:B:199:GLU:O	2:B:202:VAL:HB	0.62	1.94	19	4
1:A:34:ASN:O	1:A:37:ARG:HG2	0.62	1.94	13	13
1:A:16:ILE:HG12	1:A:19:ARG:NH2	0.62	2.10	8	1
2:B:226:HIS:O	2:B:230:LEU:HG	0.59	1.97	4	13
2:B:215:PHE:CD2	2:B:225:VAL:HG23	0.58	2.33	4	14
2:B:204:GLN:O	2:B:207:GLU:HG2	0.58	1.98	18	7
1:A:27:HIS:HB3	1:A:30:PHE:O	0.58	1.99	20	17

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:10:LYS:O	1:A:14:LYS:HG2	0.58	1.99	14	3
2:B:215:PHE:CD2	2:B:220:ILE:HG21	0.58	2.34	12	12
2:B:225:VAL:O	2:B:229:PHE:HB2	0.58	1.99	8	2
1:A:5:GLU:HB3	1:A:6:PRO:CD	0.57	2.29	7	17
1:A:18:PHE:O	1:A:22:VAL:HG23	0.57	1.98	9	15
2:B:199:GLU:OE2	2:B:228:LEU:HD21	0.57	1.99	19	1
1:A:39:LYS:HG2	2:B:207:GLU:OE1	0.56	2.00	5	2
1:A:29:GLN:NE2	2:B:217:ALA:HB3	0.56	2.16	3	1
1:A:7:SER:O	1:A:11:GLU:HG2	0.55	2.01	4	6
2:B:227:THR:O	2:B:231:LYS:HG2	0.54	2.03	16	11
2:B:205:LEU:HB3	2:B:210:VAL:HG21	0.54	1.79	3	12
1:A:23:PHE:O	1:A:37:ARG:HD2	0.53	2.03	3	2
1:A:38:LEU:O	1:A:42:ARG:HG2	0.53	2.03	14	3
1:A:35:ASN:O	1:A:39:LYS:HG2	0.53	2.03	8	2
1:A:20:ASP:HA	1:A:23:PHE:CD2	0.52	2.39	16	8
2:B:195:GLN:O	2:B:199:GLU:HG2	0.52	2.05	19	1
2:B:199:GLU:HG2	2:B:228:LEU:HD22	0.51	1.83	7	4
1:A:14:LYS:HD3	2:B:220:ILE:HG23	0.51	1.81	15	2
1:A:19:ARG:HG3	1:A:20:ASP:N	0.51	2.20	18	1
1:A:42:ARG:NE	1:A:42:ARG:HA	0.51	2.21	4	2
1:A:10:LYS:O	1:A:13:LYS:HG2	0.50	2.07	4	5
1:A:4:VAL:HG22	2:B:194:ASN:HB2	0.50	1.83	8	2
1:A:34:ASN:O	1:A:38:LEU:HG	0.50	2.05	3	6
2:B:206:PHE:CE2	2:B:234:LEU:HB3	0.50	2.41	19	1
1:A:16:ILE:O	1:A:19:ARG:HG2	0.50	2.07	12	5
1:A:39:LYS:O	1:A:43:ASP:HB2	0.49	2.07	10	1
2:B:216:ILE:HD11	2:B:222:PRO:HB3	0.49	1.84	19	2
2:B:196:LYS:N	2:B:196:LYS:HD2	0.49	2.23	18	5
1:A:38:LEU:O	1:A:42:ARG:HG3	0.49	2.08	20	6
1:A:8:VAL:HG12	1:A:12:TYR:CE2	0.49	2.43	19	2
1:A:19:ARG:NE	2:B:201:ALA:HB1	0.49	2.23	8	1
2:B:231:LYS:N	2:B:231:LYS:HE2	0.49	2.22	11	3
1:A:34:ASN:O	1:A:37:ARG:HG3	0.48	2.08	4	1
1:A:13:LYS:HG3	1:A:14:LYS:N	0.48	2.23	17	1
2:B:224:VAL:O	2:B:228:LEU:HG	0.48	2.09	7	1
2:B:221:GLU:HB2	2:B:224:VAL:HG23	0.48	1.84	6	1
1:A:37:ARG:O	1:A:41:LEU:HG	0.47	2.09	14	3
2:B:227:THR:O	2:B:231:LYS:HD2	0.47	2.08	1	2
1:A:9:GLU:O	1:A:13:LYS:HG3	0.47	2.10	19	1
2:B:211:ARG:H	2:B:214:ASP:HB2	0.47	1.70	11	6
1:A:5:GLU:HB2	1:A:6:PRO:HD3	0.47	1.86	4	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:195:GLN:O	2:B:199:GLU:HG3	0.47	2.10	7	6
1:A:43:ASP:HB2	2:B:204:GLN:NE2	0.47	2.25	14	1
2:B:196:LYS:O	2:B:199:GLU:HB2	0.46	2.10	16	2
1:A:39:LYS:HE2	1:A:43:ASP:OD1	0.46	2.10	19	1
1:A:36:VAL:HG11	2:B:208:VAL:HG12	0.46	1.88	8	1
2:B:202:VAL:HG21	2:B:228:LEU:HD13	0.46	1.88	10	3
2:B:206:PHE:CD1	2:B:234:LEU:HD13	0.46	2.46	8	1
1:A:5:GLU:CB	1:A:6:PRO:CD	0.45	2.94	17	12
2:B:212:PHE:O	2:B:216:ILE:HG13	0.45	2.11	1	1
1:A:5:GLU:HB3	1:A:6:PRO:HD2	0.45	1.88	17	1
1:A:5:GLU:OE1	1:A:8:VAL:HG23	0.45	2.10	4	1
1:A:40:PHE:HA	2:B:204:GLN:HG2	0.45	1.89	7	3
1:A:39:LYS:HG3	2:B:207:GLU:OE2	0.45	2.12	8	1
1:A:15:ILE:O	1:A:19:ARG:HG2	0.44	2.12	16	1
1:A:14:LYS:HE3	2:B:221:GLU:OE2	0.44	2.12	3	1
1:A:34:ASN:O	1:A:37:ARG:HB3	0.44	2.12	19	2
1:A:40:PHE:CE2	2:B:205:LEU:HD23	0.44	2.48	18	1
1:A:11:GLU:O	1:A:15:ILE:HB	0.44	2.13	20	1
1:A:33:PRO:HD2	1:A:36:VAL:CG2	0.44	2.42	5	1
1:A:14:LYS:HE3	2:B:220:ILE:HG23	0.44	1.90	7	1
2:B:202:VAL:HG21	2:B:228:LEU:CD1	0.43	2.42	18	3
1:A:18:PHE:CE2	2:B:218:GLU:HG3	0.43	2.48	1	2
2:B:227:THR:O	2:B:230:LEU:HB2	0.43	2.13	13	2
2:B:193:THR:OG1	2:B:196:LYS:HD3	0.43	2.14	11	2
1:A:40:PHE:HA	2:B:204:GLN:CG	0.43	2.43	5	1
1:A:10:LYS:HD2	1:A:10:LYS:C	0.43	2.38	8	1
1:A:21:THR:HG22	1:A:27:HIS:HA	0.43	1.89	11	1
1:A:10:LYS:O	1:A:14:LYS:HG3	0.43	2.14	10	1
2:B:210:VAL:O	2:B:211:ARG:HD2	0.43	2.14	11	1
1:A:39:LYS:HD3	2:B:207:GLU:OE1	0.43	2.13	1	1
1:A:19:ARG:HH11	2:B:201:ALA:HB2	0.43	1.74	4	1
2:B:198:ALA:O	2:B:202:VAL:HG23	0.42	2.14	17	1
1:A:7:SER:O	1:A:11:GLU:HG3	0.42	2.15	17	2
1:A:16:ILE:HA	1:A:19:ARG:HG2	0.42	1.92	9	1
2:B:196:LYS:O	2:B:200:LYS:HG2	0.42	2.14	5	1
1:A:16:ILE:HA	1:A:19:ARG:NE	0.42	2.30	3	1
2:B:210:VAL:HG11	2:B:215:PHE:CE1	0.41	2.50	8	1
1:A:22:VAL:O	1:A:32:VAL:HG22	0.41	2.15	20	1
1:A:43:ASP:HB2	2:B:204:GLN:CD	0.40	2.41	4	2
2:B:199:GLU:CD	2:B:228:LEU:HD21	0.40	2.41	19	1
1:A:16:ILE:HG12	1:A:19:ARG:CZ	0.40	2.45	19	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:5:GLU:HG3	2:B:195:GLN:OE1	0.40	2.16	2	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	40/55 (73%)	36±1 (91±1%)	2±1 (4±2%)	2±1 (5±2%)	3	24
2	B	42/49 (86%)	42±0 (100±0%)	0±0 (0±0%)	0±0 (0±0%)	100	100
All	All	1640/2080 (79%)	1564 (95%)	36 (2%)	40 (2%)	7	44

All 3 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	5	GLU	20
1	A	4	VAL	16
1	A	33	PRO	4

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	39/52 (75%)	38±1 (98±2%)	1±1 (2±2%)	45	90
2	B	36/40 (90%)	35±1 (97±3%)	1±1 (3±3%)	38	86
All	All	1500/1840 (82%)	1462 (97%)	38 (3%)	42	88

All 16 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)
2	B	231	LYS	6
2	B	207	GLU	5
1	A	14	LYS	4
1	A	39	LYS	4
2	B	196	LYS	3
2	B	215	PHE	3
1	A	10	LYS	3
1	A	13	LYS	2
1	A	5	GLU	1
1	A	37	ARG	1
2	B	218	GLU	1
2	B	204	GLN	1
1	A	11	GLU	1
2	B	194	ASN	1
1	A	9	GLU	1
2	B	195	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 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 [i](#)

The completeness of assignment taking into account all chemical shift lists is 88% for the well-defined parts and 86% for the entire structure.

7.1 Chemical shift list 1

File name: `working_cs.cif`

Chemical shift list name: *starch_output*

7.1.1 Bookkeeping [i](#)

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	1117
Number of shifts mapped to atoms	1117
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 [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	89	-0.45 ± 0.15	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	83	0.29 ± 0.09	None needed (< 0.5 ppm)
$^{13}\text{C}'$	82	-0.38 ± 0.16	None needed (< 0.5 ppm)
^{15}N	82	-0.11 ± 0.30	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 88%, i.e. 1052 atoms were assigned a chemical shift out of a possible 1190. 0 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	395/408 (97%)	161/165 (98%)	158/164 (96%)	76/79 (96%)
Sidechain	601/677 (89%)	410/437 (94%)	185/212 (87%)	6/28 (21%)

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	Total	¹H	¹³C	¹⁵N
Aromatic	56/105 (53%)	43/52 (83%)	13/49 (27%)	0/4 (0%)
Overall	1052/1190 (88%)	614/654 (94%)	356/425 (84%)	82/111 (74%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 86%, i.e. 1117 atoms were assigned a chemical shift out of a possible 1303. 0 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	429/465 (92%)	176/189 (93%)	171/186 (92%)	82/90 (91%)
Sidechain	632/733 (86%)	431/473 (91%)	195/229 (85%)	6/31 (19%)
Aromatic	56/105 (53%)	43/52 (83%)	13/49 (27%)	0/4 (0%)
Overall	1117/1303 (86%)	650/714 (91%)	379/464 (82%)	88/125 (70%)

7.1.4 Statistically unusual chemical shifts [i](#)

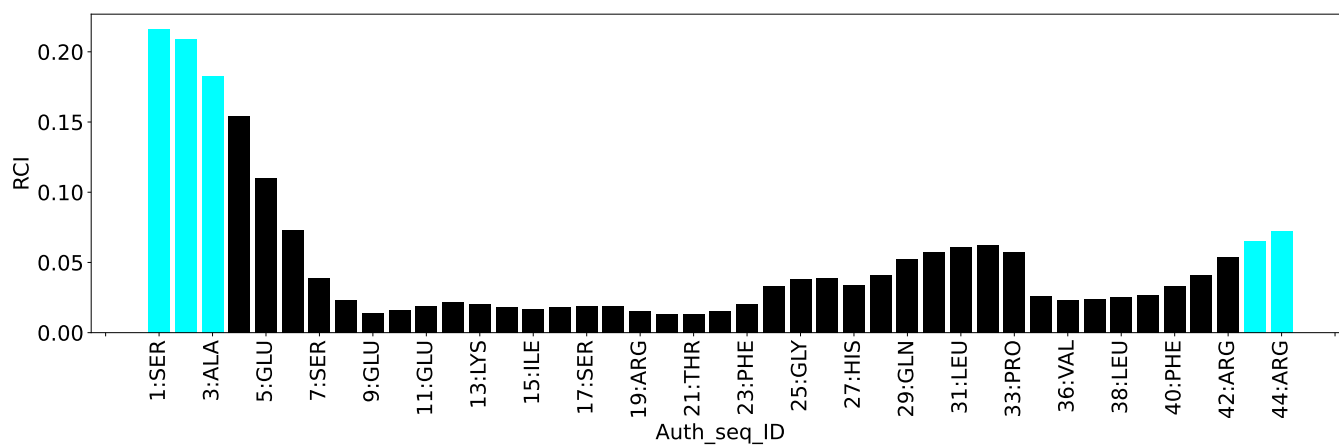
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	B	218	GLU	HG2	1.08	1.24 – 3.30	-5.8

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:



Random coil index (RCI) for chain B:

