



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 06:49 PM UTC

PDB ID : 7QB0 / pdb_00007qb0
BMRB ID : 51145
Title : Solution structure of paxillin LIM2/3
Authors : Prestel, A.; Michaelis, M.; Klishin, N.; Moeller, H.M.
Deposited on : 2021-11-17

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

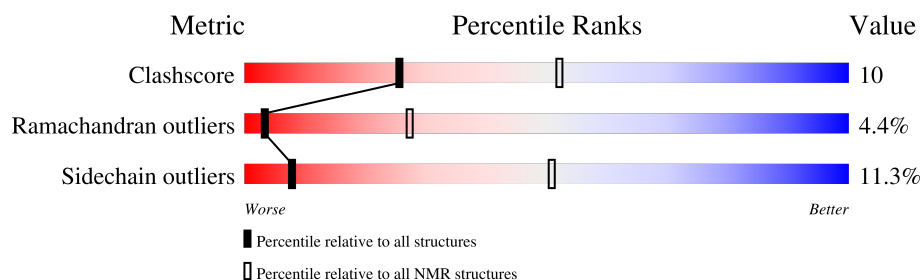
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 91%.

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	127	69% 21% • 8%

2 Ensemble composition and analysis

This entry contains 20 models. Model 3 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:1-A:117 (117)	1.90	3

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 5 clusters and 1 single-model cluster was found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1967 atoms, of which 927 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Isoform Alpha of Paxillin.

Mol	Chain	Residues	Atoms						Trace
1	A	127	Total	C	H	N	O	S	0
			1963	665	927	177	182	12	

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	121	PRO	-	expression tag	UNP P49023
A	122	GLU	-	expression tag	UNP P49023
A	123	TYR	-	expression tag	UNP P49023
A	124	PHE	-	expression tag	UNP P49023
A	125	LYS	-	expression tag	UNP P49023
A	126	ASN	-	expression tag	UNP P49023
A	127	GLY	-	expression tag	UNP P49023

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

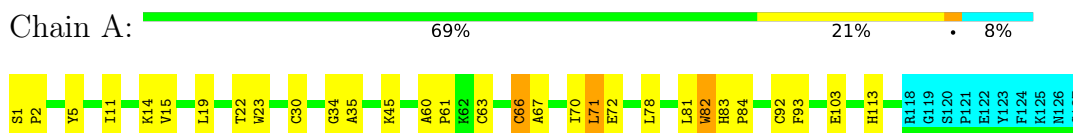
Mol	Chain	Residues	Atoms	
2	A	4	Total	Zn
			4	4

4 Residue-property plots

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: Isoform Alpha of Paxillin

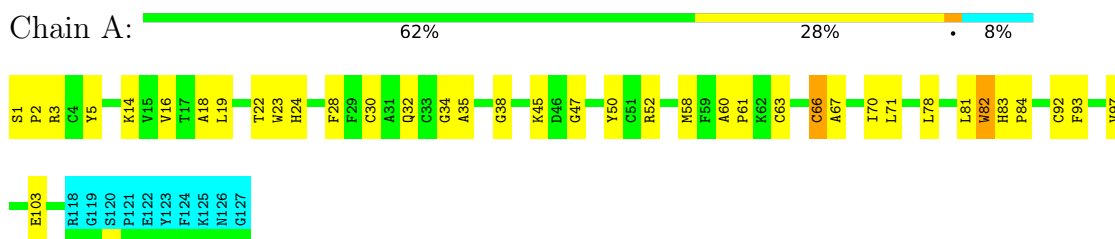


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

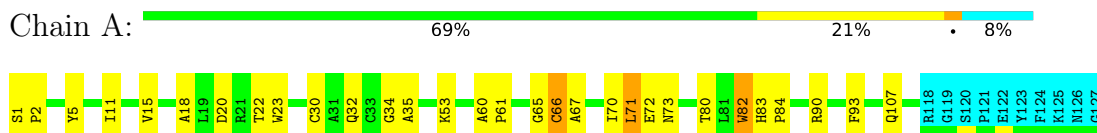
4.2.1 Score per residue for model 1

- Molecule 1: Isoform Alpha of Paxillin



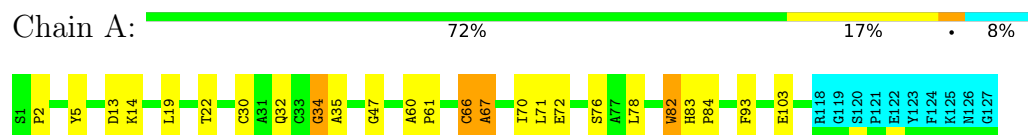
4.2.2 Score per residue for model 2

- Molecule 1: Isoform Alpha of Paxillin



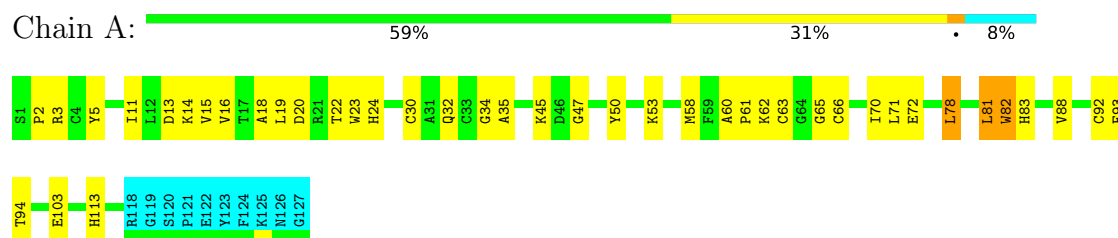
4.2.3 Score per residue for model 3 (medoid)

- Molecule 1: Isoform Alpha of Paxillin



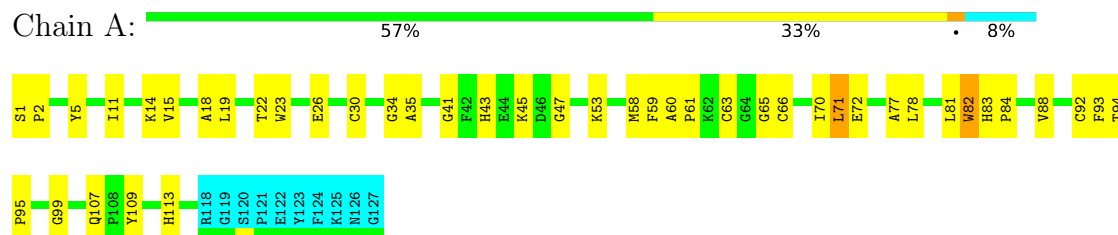
4.2.4 Score per residue for model 4

- Molecule 1: Isoform Alpha of Paxillin



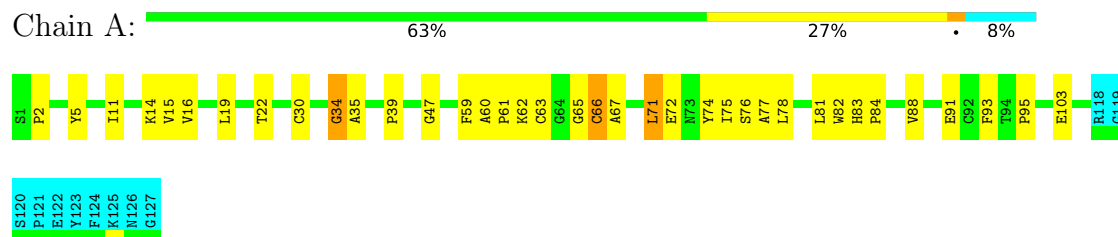
4.2.5 Score per residue for model 5

- Molecule 1: Isoform Alpha of Paxillin



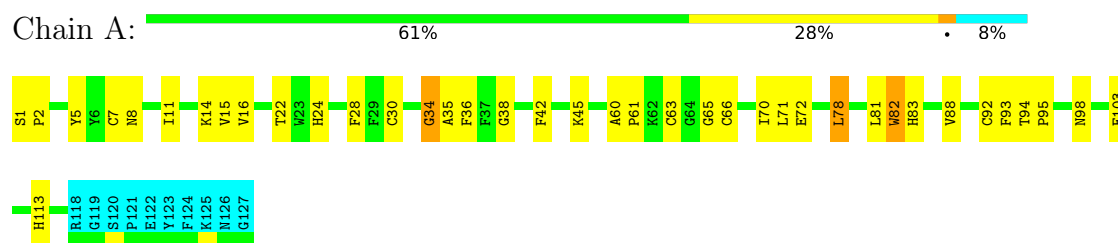
4.2.6 Score per residue for model 6

- Molecule 1: Isoform Alpha of Paxillin



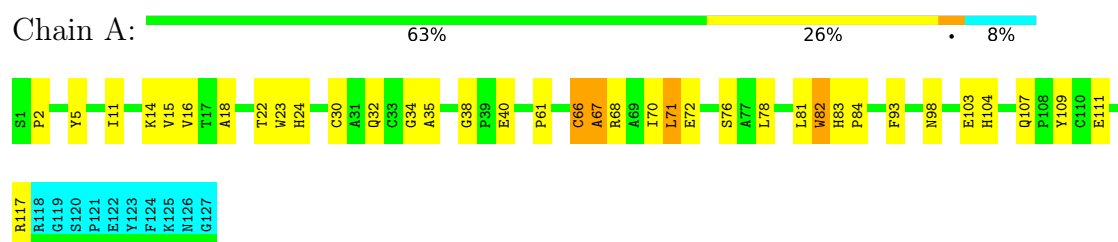
4.2.7 Score per residue for model 7

- Molecule 1: Isoform Alpha of Paxillin



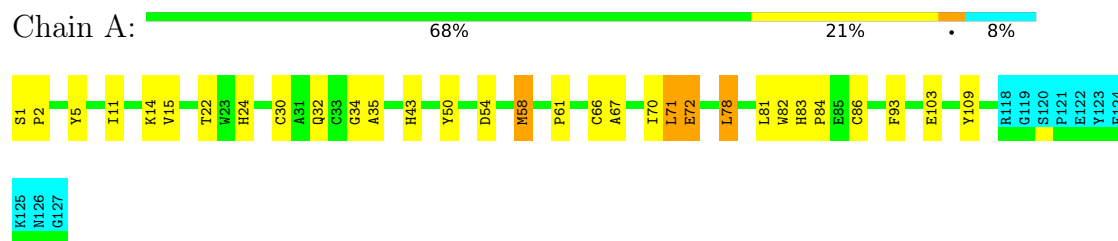
4.2.8 Score per residue for model 8

- Molecule 1: Isoform Alpha of Paxillin



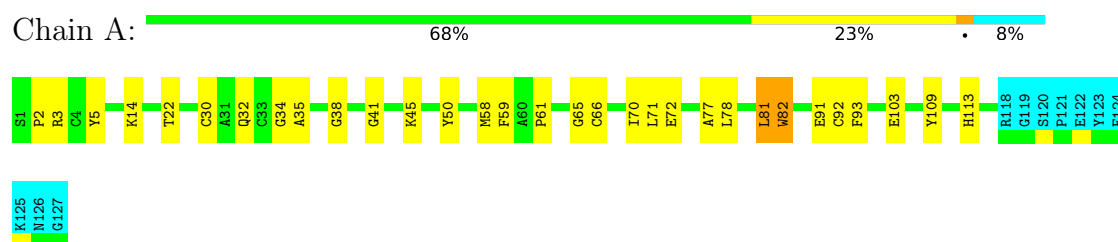
4.2.9 Score per residue for model 9

- Molecule 1: Isoform Alpha of Paxillin



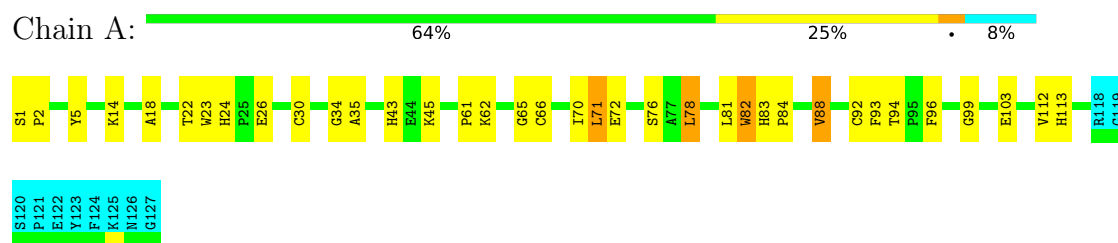
4.2.10 Score per residue for model 10

- Molecule 1: Isoform Alpha of Paxillin



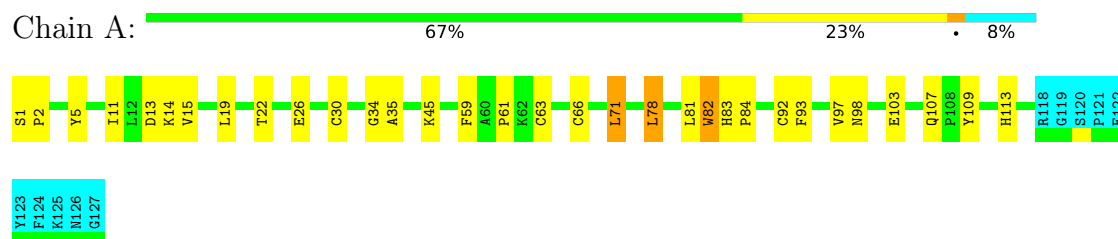
4.2.11 Score per residue for model 11

- Molecule 1: Isoform Alpha of Paxillin



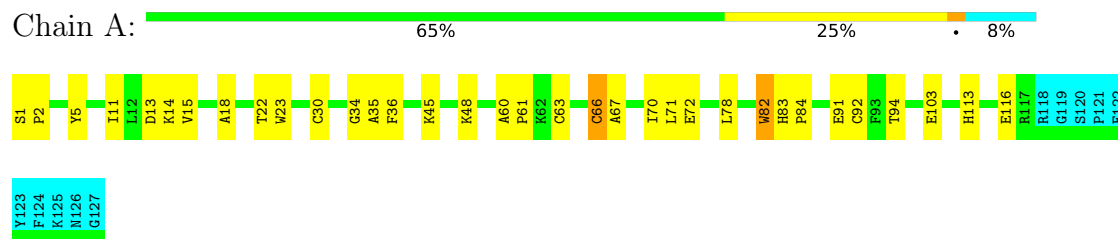
4.2.12 Score per residue for model 12

- Molecule 1: Isoform Alpha of Paxillin



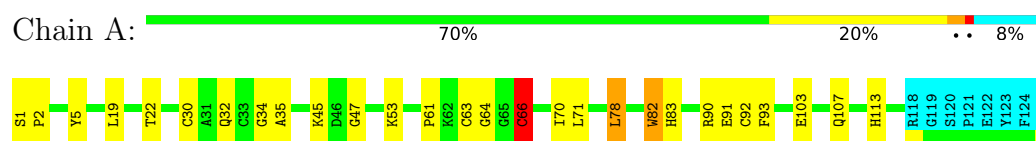
4.2.13 Score per residue for model 13

- Molecule 1: Isoform Alpha of Paxillin



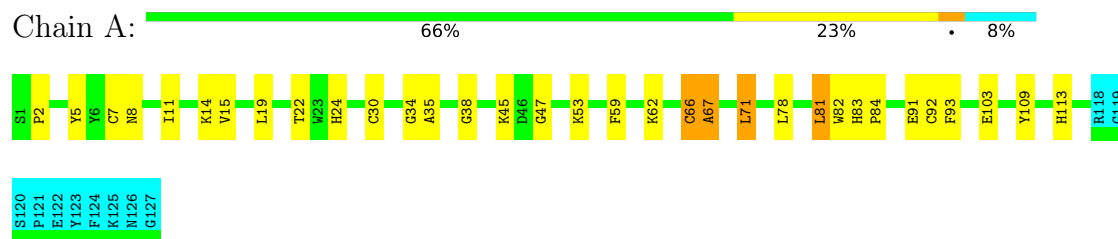
4.2.14 Score per residue for model 14

- Molecule 1: Isoform Alpha of Paxillin



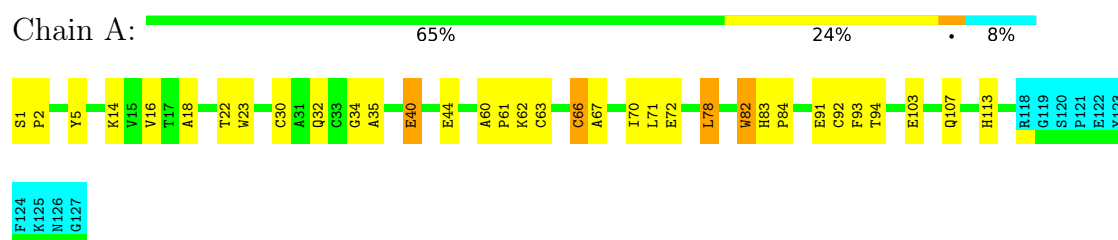
4.2.15 Score per residue for model 15

- Molecule 1: Isoform Alpha of Paxillin



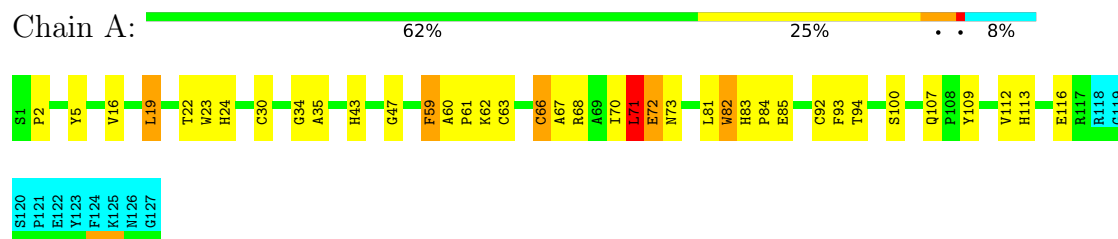
4.2.16 Score per residue for model 16

- Molecule 1: Isoform Alpha of Paxillin



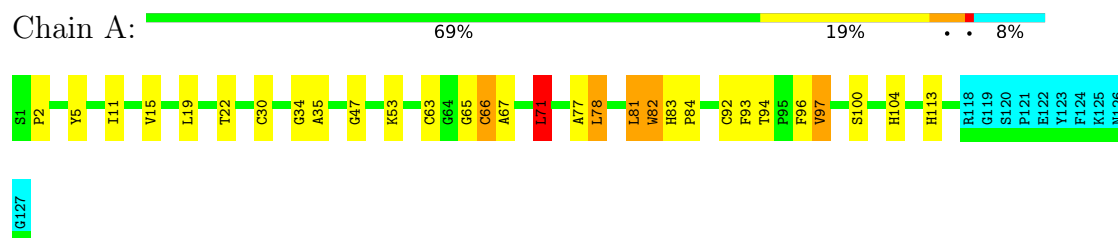
4.2.17 Score per residue for model 17

- Molecule 1: Isoform Alpha of Paxillin



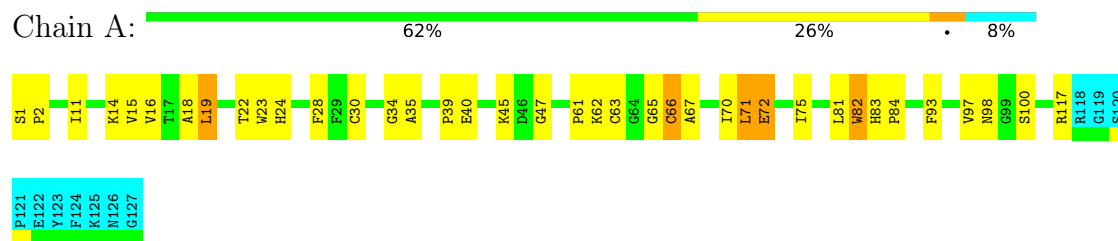
4.2.18 Score per residue for model 18

- Molecule 1: Isoform Alpha of Paxillin



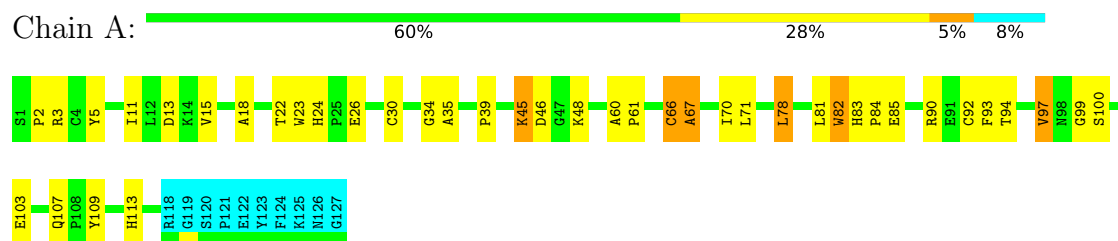
4.2.19 Score per residue for model 19

- Molecule 1: Isoform Alpha of Paxillin



4.2.20 Score per residue for model 20

- Molecule 1: Isoform Alpha of Paxillin



5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *target function*.

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

Software name	Classification	Version
CYANA	refinement	3.0
CYANA	structure calculation	3.0

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	1529
Number of shifts mapped to atoms	1529
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	91%

6 Model quality

6.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section:
ZN

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

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	955	853	855	19±3
All	All	19180	17060	17100	375

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 10.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:70:ILE:HG23	1:A:83:HIS:CD2	0.70	2.21	11	7
1:A:61:PRO:CB	1:A:81:LEU:HD12	0.69	2.18	6	1
1:A:2:PRO:HB3	1:A:22:THR:HG21	0.68	1.64	1	20
1:A:61:PRO:HB3	1:A:81:LEU:HD13	0.66	1.67	10	4
1:A:50:TYR:CE1	1:A:58:MET:HE1	0.66	2.25	1	3
1:A:66:CYS:SG	1:A:67:ALA:N	0.65	2.68	3	12
1:A:93:PHE:CG	1:A:93:PHE:O	0.64	2.50	7	13
1:A:11:ILE:CG2	1:A:15:VAL:HG22	0.63	2.24	20	4
1:A:97:VAL:HG13	1:A:100:SER:O	0.63	1.93	20	1
1:A:18:ALA:HB3	1:A:23:TRP:CD1	0.63	2.28	4	10
1:A:71:LEU:H	1:A:71:LEU:HD13	0.62	1.55	18	1
1:A:66:CYS:O	1:A:67:ALA:C	0.61	2.43	20	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:93:PHE:O	1:A:93:PHE:CG	0.61	2.53	5	6
1:A:19:LEU:HD21	1:A:47:GLY:CA	0.61	2.25	15	4
1:A:61:PRO:HB3	1:A:81:LEU:HD12	0.61	1.72	7	4
1:A:92:CYS:O	1:A:92:CYS:SG	0.60	2.59	13	1
1:A:18:ALA:HB3	1:A:23:TRP:NE1	0.60	2.12	19	10
1:A:97:VAL:HG13	1:A:100:SER:C	0.59	2.22	20	1
1:A:61:PRO:HB2	1:A:70:ILE:HD12	0.59	1.75	14	16
1:A:71:LEU:HD13	1:A:71:LEU:H	0.58	1.58	15	1
1:A:71:LEU:HD21	1:A:84:PRO:CD	0.57	2.28	19	3
1:A:19:LEU:HD21	1:A:47:GLY:HA2	0.57	1.76	14	4
1:A:11:ILE:HD13	1:A:15:VAL:HG13	0.57	1.77	13	4
1:A:30:CYS:O	1:A:34:GLY:HA2	0.56	2.00	16	20
1:A:82:TRP:CD1	1:A:82:TRP:N	0.56	2.73	2	14
1:A:2:PRO:CB	1:A:22:THR:HG21	0.56	2.30	1	19
1:A:82:TRP:N	1:A:82:TRP:CD1	0.56	2.73	18	6
1:A:74:TYR:CD2	1:A:81:LEU:HD13	0.56	2.35	6	1
1:A:78:LEU:HD22	1:A:103:GLU:CD	0.55	2.26	16	8
1:A:30:CYS:O	1:A:34:GLY:CA	0.55	2.55	19	20
1:A:83:HIS:HB3	1:A:84:PRO:HD2	0.54	1.78	12	14
1:A:60:ALA:HB1	1:A:61:PRO:HD2	0.54	1.80	7	9
1:A:19:LEU:HD11	1:A:47:GLY:O	0.54	2.02	1	3
1:A:77:ALA:C	1:A:78:LEU:HD22	0.53	2.27	18	1
1:A:16:VAL:HG13	1:A:40:GLU:HG2	0.53	1.80	16	1
1:A:19:LEU:HD21	1:A:47:GLY:O	0.53	2.04	3	1
1:A:45:LYS:O	1:A:46:ASP:CG	0.52	2.53	20	1
1:A:16:VAL:HG22	1:A:39:PRO:HG3	0.51	1.82	6	1
1:A:61:PRO:CB	1:A:81:LEU:HD13	0.51	2.36	10	1
1:A:71:LEU:HD21	1:A:84:PRO:HD2	0.51	1.82	5	3
1:A:78:LEU:HD21	1:A:103:GLU:HG3	0.50	1.82	3	6
1:A:71:LEU:CD1	1:A:71:LEU:N	0.49	2.75	17	1
1:A:92:CYS:SG	1:A:93:PHE:N	0.49	2.86	11	11
1:A:71:LEU:C	1:A:71:LEU:HD22	0.49	2.32	17	1
1:A:77:ALA:CB	1:A:78:LEU:HD22	0.49	2.38	18	1
1:A:11:ILE:HG21	1:A:15:VAL:HG22	0.49	1.84	13	4
1:A:97:VAL:HG12	1:A:99:GLY:H	0.48	1.69	20	1
1:A:112:VAL:HG22	1:A:116:GLU:CD	0.48	2.34	17	1
1:A:11:ILE:HG21	1:A:15:VAL:HB	0.48	1.85	7	1
1:A:96:PHE:O	1:A:97:VAL:C	0.47	2.57	18	1
1:A:11:ILE:HD13	1:A:15:VAL:HG23	0.47	1.86	15	8
1:A:81:LEU:N	1:A:81:LEU:CD2	0.47	2.78	4	1
1:A:71:LEU:O	1:A:72:GLU:C	0.47	2.58	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:81:LEU:N	1:A:81:LEU:HD12	0.47	2.25	19	2
1:A:61:PRO:CG	1:A:81:LEU:HD12	0.47	2.39	6	1
1:A:88:VAL:HG11	1:A:94:THR:C	0.47	2.35	7	4
1:A:54:ASP:O	1:A:58:MET:HE3	0.46	2.10	9	1
1:A:77:ALA:HB1	1:A:78:LEU:HD22	0.46	1.86	18	1
1:A:50:TYR:CD1	1:A:58:MET:HE1	0.46	2.45	9	1
1:A:71:LEU:C	1:A:71:LEU:CD2	0.45	2.89	15	2
1:A:71:LEU:HD22	1:A:73:ASN:N	0.45	2.26	17	1
1:A:88:VAL:HG11	1:A:95:PRO:N	0.45	2.26	7	2
1:A:78:LEU:HD22	1:A:103:GLU:CA	0.45	2.41	12	1
1:A:92:CYS:C	1:A:94:THR:H	0.45	2.20	20	3
1:A:19:LEU:HD22	1:A:47:GLY:HA2	0.45	1.88	17	1
1:A:63:CYS:HB3	1:A:66:CYS:O	0.45	2.12	14	1
1:A:92:CYS:HB3	1:A:113:HIS:CE1	0.44	2.47	12	13
1:A:92:CYS:C	1:A:94:THR:N	0.44	2.75	13	6
1:A:63:CYS:HB2	1:A:83:HIS:CE1	0.44	2.47	7	4
1:A:45:LYS:NZ	1:A:58:MET:HE1	0.44	2.28	5	1
1:A:7:CYS:O	1:A:8:ASN:C	0.44	2.61	15	2
1:A:70:ILE:O	1:A:71:LEU:C	0.44	2.61	8	1
1:A:77:ALA:HB1	1:A:78:LEU:HD12	0.44	1.90	10	3
1:A:81:LEU:HD12	1:A:81:LEU:N	0.44	2.28	20	1
1:A:61:PRO:HG3	1:A:81:LEU:HD12	0.43	1.89	6	1
1:A:53:LYS:CE	1:A:80:THR:HG23	0.43	2.43	2	1
1:A:71:LEU:N	1:A:71:LEU:HD22	0.43	2.27	18	1
1:A:11:ILE:HG22	1:A:15:VAL:HG22	0.43	1.90	12	3
1:A:71:LEU:HD11	1:A:73:ASN:OD1	0.42	2.14	2	1
1:A:11:ILE:CD1	1:A:15:VAL:HG13	0.42	2.42	4	1
1:A:16:VAL:HG23	1:A:23:TRP:HB2	0.42	1.90	17	3
1:A:53:LYS:NZ	1:A:80:THR:HG23	0.42	2.29	2	1
1:A:16:VAL:HG11	1:A:28:PHE:CE1	0.42	2.49	1	1
1:A:59:PHE:O	1:A:60:ALA:HB3	0.42	2.15	17	1
1:A:81:LEU:H	1:A:81:LEU:HD22	0.41	1.76	18	2
1:A:16:VAL:HG11	1:A:28:PHE:CD1	0.41	2.50	1	3
1:A:83:HIS:HB2	1:A:86:CYS:SG	0.41	2.55	9	1
1:A:81:LEU:N	1:A:81:LEU:CD1	0.41	2.83	19	2
1:A:88:VAL:HG12	1:A:95:PRO:N	0.41	2.31	6	1
1:A:19:LEU:HD22	1:A:19:LEU:N	0.41	2.31	12	1
1:A:78:LEU:HD22	1:A:103:GLU:HA	0.41	1.92	12	1
1:A:81:LEU:HD22	1:A:81:LEU:N	0.40	2.31	18	1
1:A:97:VAL:HG23	1:A:97:VAL:O	0.40	2.16	12	1
1:A:112:VAL:HG13	1:A:113:HIS:N	0.40	2.32	11	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	116/127 (91%)	99±2 (85±1%)	12±2 (11±2%)	5±1 (4±1%)	3	27
All	All	2320/2540 (91%)	1972 (85%)	246 (11%)	102 (4%)	3	27

All 19 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	35	ALA	20
1	A	71	LEU	18
1	A	72	GLU	14
1	A	65	GLY	9
1	A	38	GLY	5
1	A	13	ASP	5
1	A	67	ALA	5
1	A	66	CYS	5
1	A	97	VAL	4
1	A	34	GLY	3
1	A	100	SER	3
1	A	41	GLY	2
1	A	99	GLY	2
1	A	39	PRO	2
1	A	60	ALA	1
1	A	98	ASN	1
1	A	96	PHE	1
1	A	64	GLY	1
1	A	40	GLU	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	100/108 (93%)	89±2 (89±2%)	11±2 (11±2%)	8	51
All	All	2000/2160 (93%)	1774 (89%)	226 (11%)	8	51

All 42 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	5	TYR	19
1	A	82	TRP	17
1	A	66	CYS	16
1	A	14	LYS	15
1	A	1	SER	11
1	A	45	LYS	11
1	A	24	HIS	10
1	A	32	GLN	9
1	A	78	LEU	9
1	A	63	CYS	8
1	A	107	GLN	8
1	A	109	TYR	8
1	A	62	LYS	7
1	A	59	PHE	6
1	A	91	GLU	6
1	A	53	LYS	5
1	A	81	LEU	5
1	A	71	LEU	5
1	A	3	ARG	4
1	A	76	SER	4
1	A	26	GLU	4
1	A	43	HIS	4
1	A	90	ARG	3
1	A	98	ASN	3
1	A	20	ASP	2
1	A	75	ILE	2
1	A	36	PHE	2
1	A	40	GLU	2
1	A	68	ARG	2
1	A	104	HIS	2
1	A	117	ARG	2
1	A	72	GLU	2
1	A	48	LYS	2
1	A	19	LEU	2
1	A	85	GLU	2
1	A	52	ARG	1

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Mol	Chain	Res	Type	Models (Total)
1	A	42	PHE	1
1	A	111	GLU	1
1	A	58	MET	1
1	A	88	VAL	1
1	A	116	GLU	1
1	A	44	GLU	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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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 91% for the well-defined parts and 90% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chemical_shifts_1*

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	1529
Number of shifts mapped to atoms	1529
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	3

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
$^{13}\text{C}_\alpha$	127	0.08 ± 0.15	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	116	0.16 ± 0.19	None needed (< 0.5 ppm)
$^{13}\text{C}'$	125	0.28 ± 0.17	None needed (< 0.5 ppm)
^{15}N	115	0.26 ± 0.38	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 91%, i.e. 1412 atoms were assigned a chemical shift out of a possible 1560. 0 out of 10 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	572/578 (99%)	233/235 (99%)	232/234 (99%)	107/109 (98%)
Sidechain	638/726 (88%)	433/468 (93%)	199/229 (87%)	6/29 (21%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	202/256 (79%)	101/128 (79%)	99/119 (83%)	2/9 (22%)
Overall	1412/1560 (91%)	767/831 (92%)	530/582 (91%)	115/147 (78%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 90%, i.e. 1529 atoms were assigned a chemical shift out of a possible 1692. 0 out of 10 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	620/628 (99%)	253/256 (99%)	252/254 (99%)	115/118 (97%)
Sidechain	689/789 (87%)	467/507 (92%)	215/248 (87%)	7/34 (21%)
Aromatic	220/275 (80%)	110/137 (80%)	108/129 (84%)	2/9 (22%)
Overall	1529/1692 (90%)	830/900 (92%)	575/631 (91%)	124/161 (77%)

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	A	42	PHE	HB3	0.52	1.03 – 4.85	-6.3
1	A	101	PHE	HB3	0.59	1.03 – 4.85	-6.2
1	A	64	GLY	HA3	2.08	2.08 – 5.71	-5.0

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:

