



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 01:27 AM UTC

PDB ID : 2M32 / pdb_00002m32
BMRB ID : 18942
Title : Alpha-1 integrin I-domain in complex with GLOGEN triple helical peptide
Authors : Chin, Y.; Headey, S.; Mohanty, B.; McEwan, P.; Swarbrick, J.; Mulhern, T.;
Emsley, J.; Simpson, J.; Scanlon, M.
Deposited on : 2013-01-07

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
Mogul : 2022.3.0, CSD as543be (2022)
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

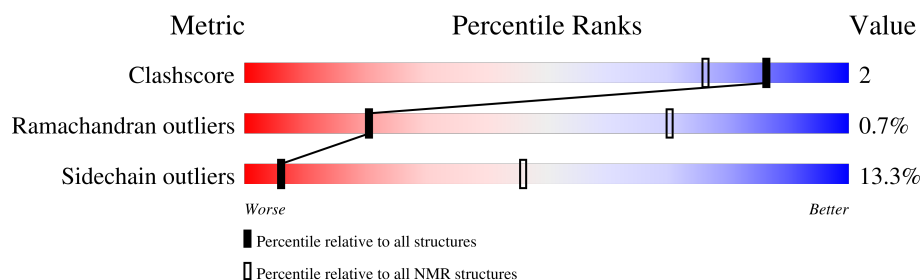
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 70%.

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	192	83% 11% 5%
2	B	23	65% 35%
2	C	23	65% 35%
2	D	23	65% 35%

2 Ensemble composition and analysis

This entry contains 10 models. Model 8 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:2-A:144, A:154-A:192, B:102-B:103, B:105-B:106, B:108-B:109, B:111-B:115, B:117-B:118, B:120-B:121, C:202-C:203, C:205-C:206, C:208-C:209, C:211-C:215, C:217-C:218, C:220-C:221, D:302-D:303, D:305-D:306, D:308-D:309, D:311-D:315, D:317-D:318, D:320-D:321 (227)	1.16	8

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 1 single-model cluster was found.

Cluster number	Models
1	2, 3, 5, 8, 10
2	1, 6, 7, 9
Single-model clusters	4

3 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3837 atoms, of which 1905 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Integrin alpha-1.

Mol	Chain	Residues	Atoms						Trace
1	A	192	Total	C	H	N	O	S	0
			3038	944	1527	265	298	4	

- Molecule 2 is a protein called GLOGEN peptide.

Mol	Chain	Residues	Atoms					Trace
2	B	23	Total	C	H	N	O	1
			266	86	126	23	31	
2	C	23	Total	C	H	N	O	1
			266	86	126	23	31	
2	D	23	Total	C	H	N	O	1
			266	86	126	23	31	

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

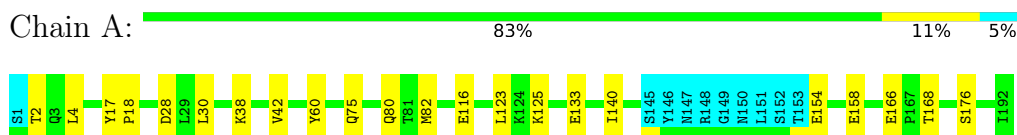
Mol	Chain	Residues	Atoms	
3	A	1	Total	Mg
			1	1

4 Residue-property plots [i](#)

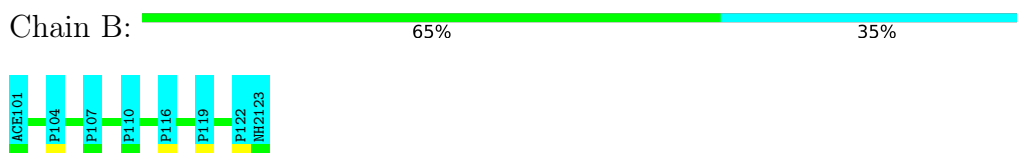
4.1 Average score per residue in the NMR ensemble

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.

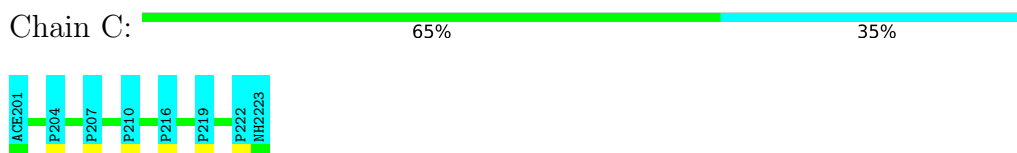
- Molecule 1: Integrin alpha-1



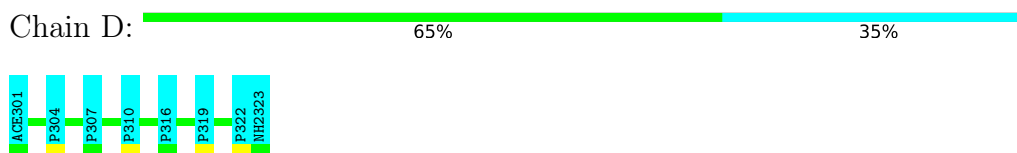
- Molecule 2: GLOGEN peptide



- Molecule 2: GLOGEN peptide



- Molecule 2: GLOGEN peptide



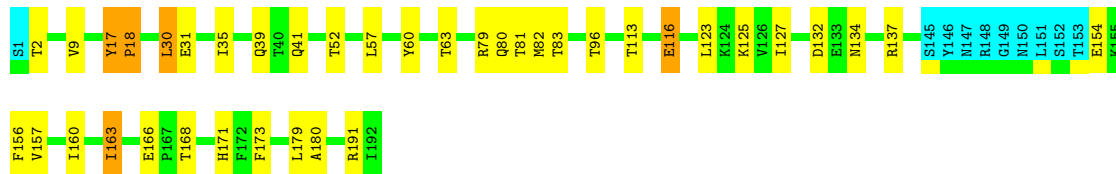
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

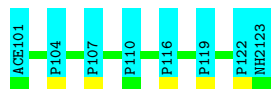
- Molecule 1: Integrin alpha-1

Chain A:  74% 18% • 5%



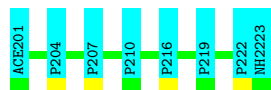
- Molecule 2: GLOGEN peptide

Chain B:  65% 35%



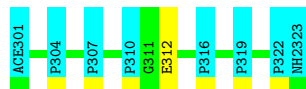
- Molecule 2: GLOGEN peptide

Chain C:  65% 35%



- Molecule 2: GLOGEN peptide

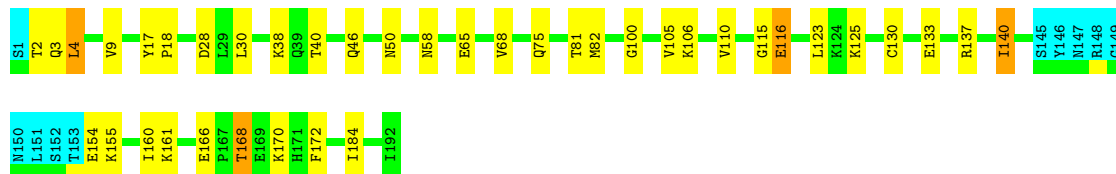
Chain D:  61% 35%



4.2.2 Score per residue for model 2

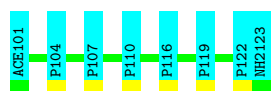
- Molecule 1: Integrin alpha-1

Chain A:  74% 18% • 5%

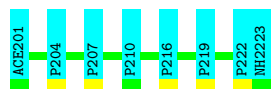


- Molecule 2: GLOGEN peptide

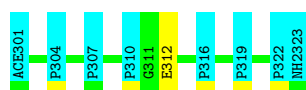
Chain B:  65% 35%



- Molecule 2: GLOGEN peptide

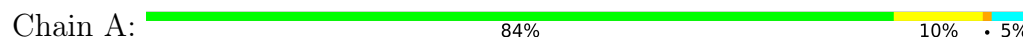


- Molecule 2: GLOGEN peptide

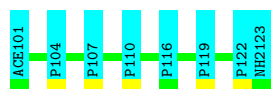


4.2.3 Score per residue for model 3

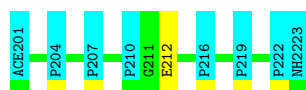
- Molecule 1: Integrin alpha-1



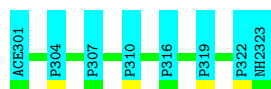
- Molecule 2: GLOGEN peptide



- Molecule 2: GLOGEN peptide

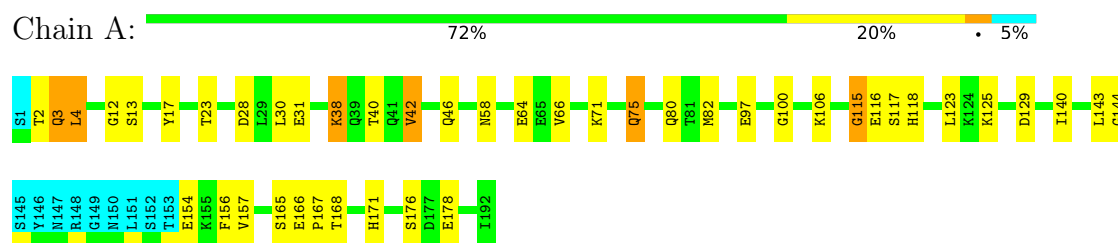


- Molecule 2: GLOGEN peptide



4.2.4 Score per residue for model 4

- Molecule 1: Integrin alpha-1



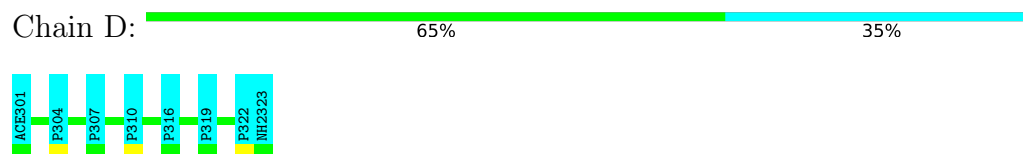
- Molecule 2: GLOGEN peptide



- Molecule 2: GLOGEN peptide

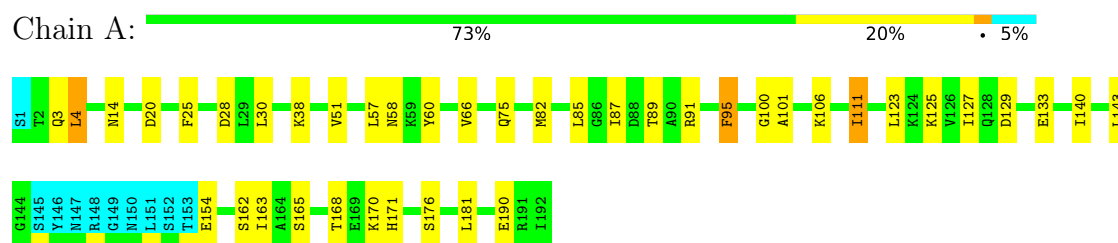


- Molecule 2: GLOGEN peptide



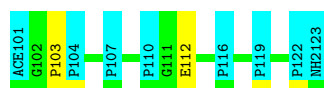
4.2.5 Score per residue for model 5

- Molecule 1: Integrin alpha-1

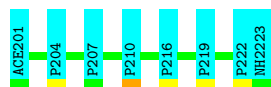


- Molecule 2: GLOGEN peptide

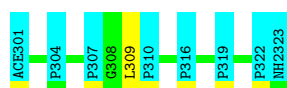




- Molecule 2: GLOGEN peptide

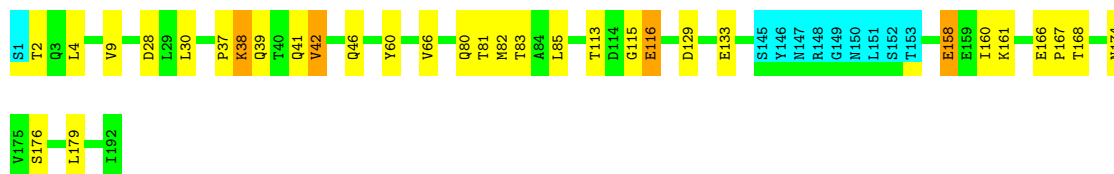
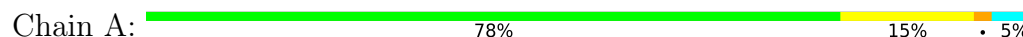


- Molecule 2: GLOGEN peptide

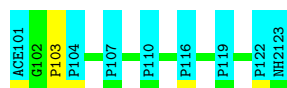


4.2.6 Score per residue for model 6

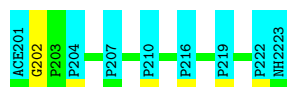
- Molecule 1: Integrin alpha-1



- Molecule 2: GLOGEN peptide

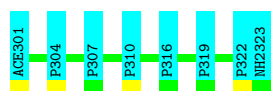


- Molecule 2: GLOGEN peptide



- Molecule 2: GLOGEN peptide

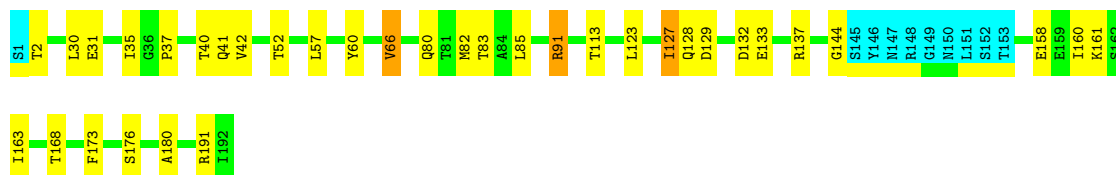




4.2.7 Score per residue for model 7

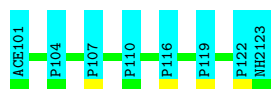
- Molecule 1: Integrin alpha-1

Chain A: 77% 17% • 5%



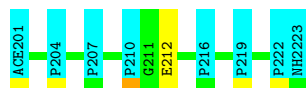
- Molecule 2: GLOGEN peptide

Chain B: 65% 35%



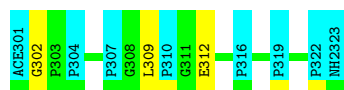
- Molecule 2: GLOGEN peptide

Chain C: 61% • 35%



- Molecule 2: GLOGEN peptide

Chain D: 52% 13% 35%



4.2.8 Score per residue for model 8 (medoid)

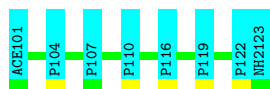
- Molecule 1: Integrin alpha-1

Chain A: 75% 18% • 5%

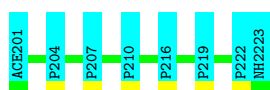




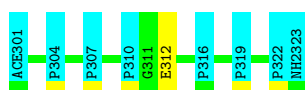
- Molecule 2: GLOGEN peptide



- Molecule 2: GLOGEN peptide

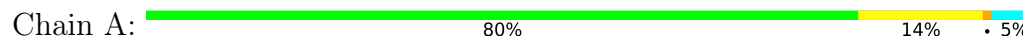


- Molecule 2: GLOGEN peptide

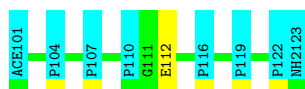


4.2.9 Score per residue for model 9

- Molecule 1: Integrin alpha-1

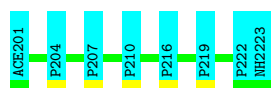


- Molecule 2: GLOGEN peptide

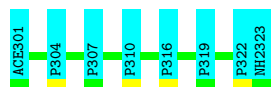


- Molecule 2: GLOGEN peptide



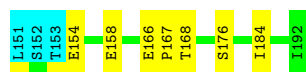
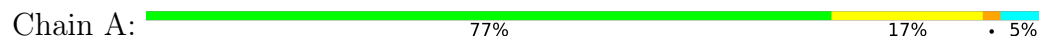


- Molecule 2: GLOGEN peptide

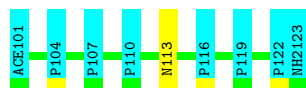


4.2.10 Score per residue for model 10

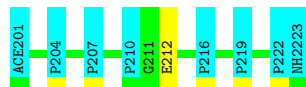
- Molecule 1: Integrin alpha-1



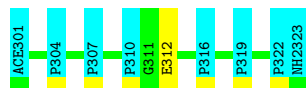
- Molecule 2: GLOGEN peptide



- Molecule 2: GLOGEN peptide



- Molecule 2: GLOGEN peptide



5 Refinement protocol and experimental data overview

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

Of the 200 calculated structures, 10 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
CNS	structure solution	
CNS	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	2148
Number of shifts mapped to atoms	2148
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	70%

6 Model quality (i)

6.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: HYP, MG, ACE, NH2

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	0.58±0.11	0±0/1453 (0.0± 0.0%)	1.09±0.16	2±4/1958 (0.1± 0.2%)
2	B	0.50±0.08	0±0/87 (0.0± 0.0%)	0.73±0.11	0±0/111 (0.0± 0.0%)
2	C	0.50±0.08	0±0/87 (0.0± 0.0%)	0.98±0.10	0±1/111 (0.2± 0.5%)
2	D	0.49±0.10	0±0/87 (0.0± 0.0%)	0.76±0.14	0±0/111 (0.0± 0.0%)
All	All	0.58	1/17140 (0.0%)	1.07	19/22910 (0.1%)

All unique bond outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	114	ASP	CG-OD1	5.75	1.36	1.25	8	1

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	77	GLY	CA-C-N	7.07	127.06	120.34	3	1
1	A	77	GLY	C-N-CA	7.07	127.06	120.34	3	1
1	A	166	GLU	CB-CA-C	5.83	121.66	110.17	3	1
1	A	156	PHE	CA-CB-CG	5.75	119.55	113.80	3	1
1	A	19	TRP	CA-C-N	5.67	128.15	120.38	3	1
1	A	19	TRP	C-N-CA	5.67	128.15	120.38	3	1
1	A	17	TYR	CA-C-N	5.48	126.69	119.84	1	1
1	A	17	TYR	C-N-CA	5.48	126.69	119.84	1	1
1	A	14	ASN	N-CA-C	-5.29	107.00	113.50	10	1
1	A	20	ASP	CA-CB-CG	5.22	117.82	112.60	9	1
1	A	96	THR	CA-C-N	5.20	127.77	120.28	3	1
1	A	96	THR	C-N-CA	5.20	127.77	120.28	3	1
1	A	12	GLY	CA-C-N	5.18	131.44	121.54	3	1
1	A	12	GLY	C-N-CA	5.18	131.44	121.54	3	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	85	LEU	CA-C-N	5.17	125.67	119.94	3	1
1	A	85	LEU	C-N-CA	5.17	125.67	119.94	3	1
2	C	212	GLU	CA-C-N	5.08	127.99	120.82	3	1
2	C	212	GLU	C-N-CA	5.08	127.99	120.82	3	1
1	A	129	ASP	CA-CB-CG	5.05	117.65	112.60	3	1

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	1435	1455	1448	7±3
2	C	88	79	79	0±0
2	B	88	79	79	0±0
2	D	88	79	79	0±0
All	All	17000	16920	16851	71

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 2.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:58:ASN:HB2	1:A:100:GLY:HA3	0.62	1.71	4	5
1:A:17:TYR:HB3	1:A:18:PRO:HD2	0.61	1.71	2	4
1:A:161:LYS:HG3	1:A:168:THR:HG21	0.56	1.76	2	1
1:A:60:TYR:HB2	1:A:66:VAL:HG22	0.52	1.81	5	1
1:A:42:VAL:HG23	1:A:66:VAL:HG11	0.51	1.81	4	3
1:A:113:THR:HG21	1:A:160:ILE:HG21	0.51	1.81	7	4
1:A:115:GLY:HA3	1:A:156:PHE:HB3	0.51	1.83	4	1
1:A:17:TYR:CZ	1:A:144:GLY:HA3	0.49	2.42	4	1
1:A:166:GLU:HB2	1:A:167:PRO:HD2	0.49	1.85	10	2
1:A:81:THR:HB	1:A:116:GLU:HG3	0.48	1.83	1	1
1:A:173:PHE:HB3	1:A:180:ALA:HB1	0.48	1.84	1	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:91:ARG:HA	1:A:95:PHE:HB2	0.48	1.86	5	2
1:A:81:THR:HB	1:A:116:GLU:HB3	0.48	1.84	2	2
1:A:60:TYR:HB3	1:A:65:GLU:HB2	0.48	1.85	9	1
1:A:4:LEU:HB3	1:A:106:LYS:HB2	0.48	1.86	2	5
1:A:91:ARG:HH12	1:A:133:GLU:HG3	0.47	1.68	7	1
1:A:25:PHE:CE1	1:A:181:LEU:HD21	0.47	2.44	8	2
1:A:60:TYR:HB3	1:A:66:VAL:HG22	0.47	1.86	7	1
1:A:123:LEU:HD13	1:A:163:ILE:HD11	0.46	1.86	1	1
1:A:168:THR:HG22	1:A:172:PHE:HB3	0.46	1.86	2	1
1:A:130:CYS:HB3	1:A:137:ARG:HD2	0.46	1.88	2	1
1:A:30:LEU:HD12	1:A:35:ILE:HD11	0.45	1.89	1	1
1:A:81:THR:CG2	1:A:116:GLU:HG2	0.45	2.42	6	1
1:A:158:GLU:HA	1:A:161:LYS:HG2	0.45	1.89	8	3
1:A:35:ILE:HD12	1:A:63:THR:HA	0.44	1.90	10	2
1:A:58:ASN:HB2	1:A:100:GLY:CA	0.44	2.42	10	1
1:A:83:THR:HG21	1:A:113:THR:HB	0.44	1.89	7	1
1:A:115:GLY:HA2	1:A:160:ILE:HD11	0.43	1.90	6	1
1:A:23:THR:HG22	1:A:71:LYS:HA	0.43	1.89	4	1
1:A:46:GLN:HG3	1:A:52:THR:HB	0.43	1.91	9	1
1:A:80:GLN:HE21	2:C:212:GLU:HB3	0.43	1.74	4	1
1:A:87:ILE:HG12	1:A:111:ILE:HG12	0.43	1.90	5	1
1:A:3:GLN:HA	1:A:3:GLN:HE21	0.42	1.74	4	1
1:A:57:LEU:HD23	1:A:101:ALA:HA	0.42	1.92	5	1
1:A:17:TYR:CB	1:A:18:PRO:HD2	0.42	2.45	1	1
1:A:12:GLY:HA3	1:A:75:GLN:HG3	0.42	1.91	4	2
1:A:110:VAL:HG13	1:A:140:ILE:HD11	0.41	1.92	2	1
1:A:166:GLU:HB3	1:A:167:PRO:HD2	0.41	1.92	6	1
1:A:166:GLU:HB2	1:A:171:HIS:CD2	0.41	2.51	1	1
1:A:168:THR:HG22	1:A:172:PHE:CB	0.41	2.46	2	1
1:A:51:VAL:HG21	1:A:89:THR:HG21	0.41	1.92	5	1
1:A:127:ILE:HG12	1:A:163:ILE:HG13	0.40	1.92	7	1
1:A:127:ILE:HG12	1:A:163:ILE:HA	0.40	1.93	5	1
1:A:59:LYS:HE2	1:A:59:LYS:HA	0.40	1.92	8	1
1:A:65:GLU:HA	1:A:68:VAL:HG12	0.40	1.92	2	1

6.3 Torsion angles ⓘ

6.3.1 Protein backbone ⓘ

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	181/192 (94%)	164±3 (91±2%)	15±3 (8±2%)	2±1 (1±0%)	18	68
2	B	15/23 (65%)	15±1 (99±4%)	0±1 (1±4%)	0±0 (0±0%)	100	100
2	C	15/23 (65%)	14±1 (97±5%)	0±1 (3±5%)	0±0 (0±0%)	100	100
2	D	15/23 (65%)	14±0 (96±3%)	1±0 (4±3%)	0±0 (0±0%)	100	100
All	All	2260/2610 (87%)	2081 (92%)	164 (7%)	15 (1%)	20	70

All 8 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	115	GLY	3
1	A	38	LYS	3
1	A	37	PRO	3
1	A	18	PRO	2
1	A	13	SER	1
1	A	134	ASN	1
1	A	144	GLY	1
1	A	116	GLU	1

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	159/168 (95%)	136±5 (85±3%)	23±5 (15±3%)	5	43
2	B	8/8 (100%)	8±0 (95±6%)	0±0 (5±6%)	23	74
2	C	8/8 (100%)	8±0 (98±5%)	0±0 (2±5%)	42	88
2	D	8/8 (100%)	8±0 (94±6%)	0±0 (6±6%)	18	67
All	All	1830/1920 (95%)	1586 (87%)	244 (13%)	6	46

All 86 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	30	LEU	9
1	A	82	MET	9
1	A	168	THR	8
1	A	2	THR	7
1	A	176	SER	7
1	A	116	GLU	6
1	A	125	LYS	6
1	A	154	GLU	6
1	A	4	LEU	6
1	A	28	ASP	6
1	A	9	VAL	5
1	A	80	GLN	5
2	D	312	GLU	5
1	A	38	LYS	5
1	A	75	GLN	5
1	A	123	LEU	5
1	A	133	GLU	5
1	A	140	ILE	5
1	A	41	GLN	4
1	A	83	THR	4
1	A	3	GLN	4
1	A	40	THR	4
1	A	129	ASP	4
1	A	31	GLU	3
1	A	39	GLN	3
1	A	52	THR	3
1	A	60	TYR	3
1	A	132	ASP	3
1	A	137	ARG	3
1	A	46	GLN	3
1	A	95	PHE	3
1	A	158	GLU	3
1	A	42	VAL	3
1	A	165	SER	3
1	A	171	HIS	3
2	B	112	GLU	3
1	A	85	LEU	3
1	A	57	LEU	2
1	A	127	ILE	2
1	A	156	PHE	2
1	A	157	VAL	2
1	A	163	ILE	2
1	A	179	LEU	2

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Mol	Chain	Res	Type	Models (Total)
1	A	191	ARG	2
1	A	166	GLU	2
1	A	170	LYS	2
1	A	184	ILE	2
1	A	17	TYR	2
1	A	118	HIS	2
1	A	128	GLN	2
1	A	13	SER	2
1	A	64	GLU	2
1	A	97	GLU	2
1	A	117	SER	2
1	A	143	LEU	2
1	A	20	ASP	2
1	A	111	ILE	2
1	A	190	GLU	2
2	C	212	GLU	2
1	A	34	ASP	2
1	A	79	ARG	1
1	A	96	THR	1
1	A	134	ASN	1
1	A	50	ASN	1
1	A	105	VAL	1
1	A	155	LYS	1
1	A	160	ILE	1
1	A	178	GLU	1
1	A	14	ASN	1
1	A	162	SER	1
1	A	174	ASN	1
1	A	35	ILE	1
1	A	66	VAL	1
1	A	91	ARG	1
1	A	11	ASP	1
1	A	55	PHE	1
1	A	59	LYS	1
1	A	63	THR	1
1	A	114	ASP	1
1	A	61	SER	1
1	A	87	ILE	1
1	A	159	GLU	1
1	A	21	SER	1
1	A	54	GLU	1
1	A	131	GLU	1

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Mol	Chain	Res	Type	Models (Total)
2	B	113	ASN	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

18 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length is the number of standard deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
2	HYP	B	110	2	7,8,9	0.80±0.12	0±0 (0±0%)
2	HYP	C	216	2	7,8,9	0.66±0.08	0±0 (0±0%)
2	HYP	D	310	2	7,8,9	0.68±0.07	0±0 (0±0%)
2	HYP	D	322	2	7,8,9	1.42±0.60	1±0 (8±6%)
2	HYP	D	304	2	7,8,9	0.62±0.06	0±0 (0±0%)
2	HYP	C	207	2	7,8,9	0.63±0.05	0±0 (0±0%)
2	HYP	D	307	2	7,8,9	0.68±0.07	0±0 (0±0%)
2	HYP	C	210	2	7,8,9	0.70±0.09	0±0 (0±0%)
2	HYP	B	119	2	7,8,9	0.61±0.06	0±0 (0±0%)
2	HYP	B	104	2	7,8,9	0.64±0.08	0±0 (0±0%)
2	HYP	B	122	2	7,8,9	1.43±0.60	1±0 (8±6%)
2	HYP	B	116	2	7,8,9	0.66±0.04	0±0 (0±0%)
2	HYP	C	222	2	7,8,9	1.42±0.59	1±0 (8±6%)
2	HYP	D	319	2	7,8,9	0.62±0.06	0±0 (0±0%)
2	HYP	B	107	2	7,8,9	0.59±0.05	0±0 (0±0%)
2	HYP	C	204	2	7,8,9	0.65±0.07	0±0 (0±0%)
2	HYP	C	219	2	7,8,9	0.61±0.05	0±0 (0±0%)
2	HYP	D	316	2	7,8,9	0.63±0.04	0±0 (0±0%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types,

if any, to which the group is linked. The Z score for a bond angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Counts	Bond angles	
						RMSZ	#Z>2
2	HYP	B	110	2	5,10,12	1.35±0.49	1±1 (12±18%)
2	HYP	C	216	2	5,10,12	1.75±0.26	1±1 (28±13%)
2	HYP	D	310	2	5,10,12	1.53±0.14	1±0 (20±0%)
2	HYP	D	322	2	5,10,12	1.77±0.18	2±0 (30±10%)
2	HYP	D	304	2	5,10,12	1.50±0.10	1±1 (18±10%)
2	HYP	C	207	2	5,10,12	1.55±0.20	1±1 (16±12%)
2	HYP	D	307	2	5,10,12	1.31±0.20	0±0 (8±9%)
2	HYP	C	210	2	5,10,12	1.38±0.22	1±1 (12±13%)
2	HYP	B	119	2	5,10,12	1.47±0.11	1±1 (18±10%)
2	HYP	B	104	2	5,10,12	1.55±0.14	1±1 (26±15%)
2	HYP	B	122	2	5,10,12	1.98±0.13	2±1 (40±12%)
2	HYP	B	116	2	5,10,12	1.51±0.17	1±0 (14±9%)
2	HYP	C	222	2	5,10,12	1.74±0.21	1±1 (28±18%)
2	HYP	D	319	2	5,10,12	1.64±0.23	1±1 (18±14%)
2	HYP	B	107	2	5,10,12	1.38±0.14	1±1 (12±13%)
2	HYP	C	204	2	5,10,12	1.57±0.14	2±0 (30±10%)
2	HYP	C	219	2	5,10,12	1.61±0.22	1±0 (20±8%)
2	HYP	D	316	2	5,10,12	1.35±0.17	1±1 (12±13%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HYP	C	207	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	C	216	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	C	222	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	C	219	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	B	119	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	D	319	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	D	307	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	C	204	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	C	210	2	-	0±0,0,11,13	0±0,1,1,1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HYP	B	104	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	B	110	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	B	107	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	B	116	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	D	316	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	B	122	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	D	322	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	D	304	2	-	0±0,0,11,13	0±0,1,1,1
2	HYP	D	310	2	-	0±0,0,11,13	0±0,1,1,1

All unique bond 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(Å)	Ideal(Å)	Models	
								Worst	Total
2	B	122	HYP	O-C	4.99	1.38	1.20	3	6
2	D	322	HYP	O-C	4.99	1.38	1.20	3	6
2	C	222	HYP	O-C	4.97	1.38	1.20	3	6

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
2	B	110	HYP	CB-CG-CD	4.15	107.78	103.16	3	2
2	D	322	HYP	O-C-CA	4.09	114.26	124.77	3	6
2	B	122	HYP	O-C-CA	4.07	114.31	124.77	3	6
2	C	222	HYP	O-C-CA	3.87	114.81	124.77	3	6
2	B	122	HYP	CB-CG-CD	3.61	107.18	103.16	1	8
2	C	219	HYP	CB-CG-CD	3.58	107.14	103.16	3	9
2	C	216	HYP	CB-CG-CD	3.54	107.10	103.16	3	8
2	D	319	HYP	CB-CG-CD	3.49	107.05	103.16	3	6
2	D	322	HYP	CB-CG-CD	3.26	106.79	103.16	2	5
2	C	207	HYP	CB-CG-CD	3.07	106.57	103.16	3	6
2	D	310	HYP	CB-CG-CD	2.97	106.46	103.16	1	7
2	C	222	HYP	CB-CG-CD	2.94	106.43	103.16	1	3
2	C	204	HYP	CB-CG-CD	2.89	106.38	103.16	8	10
2	C	216	HYP	CG-CB-CA	2.89	107.09	103.75	3	2
2	C	219	HYP	CG-CB-CA	2.86	107.06	103.75	3	1
2	D	310	HYP	OD1-CG-CD	2.82	104.54	110.30	7	3
2	B	110	HYP	CG-CB-CA	2.80	107.00	103.75	2	3
2	D	319	HYP	CG-CB-CA	2.80	106.99	103.75	3	1
2	B	116	HYP	CB-CG-CD	2.78	106.25	103.16	10	7

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
2	B	119	HYP	CB-CG-CD	2.77	106.25	103.16	3	8
2	B	104	HYP	CB-CG-CD	2.72	106.18	103.16	10	8
2	B	110	HYP	OD1-CG-CD	2.70	104.78	110.30	8	1
2	C	222	HYP	OD1-CG-CD	2.68	104.82	110.30	2	3
2	B	122	HYP	CG-CB-CA	2.66	106.83	103.75	8	4
2	C	222	HYP	CG-CB-CA	2.66	106.83	103.75	1	2
2	B	107	HYP	CB-CG-CD	2.66	106.11	103.16	3	5
2	D	322	HYP	CG-CB-CA	2.61	106.78	103.75	2	1
2	D	304	HYP	CB-CG-CD	2.60	106.05	103.16	10	8
2	C	210	HYP	CB-CG-CD	2.57	106.02	103.16	9	2
2	B	122	HYP	OD1-CG-CD	2.53	105.12	110.30	2	2
2	B	104	HYP	OD1-CG-CD	2.50	105.20	110.30	6	4
2	D	307	HYP	CB-CG-CD	2.47	105.91	103.16	5	4
2	C	207	HYP	CG-CB-CA	2.43	106.56	103.75	3	1
2	D	316	HYP	CB-CG-CD	2.42	105.85	103.16	10	4
2	C	204	HYP	OD1-CG-CD	2.35	105.50	110.30	10	4
2	C	216	HYP	OD1-CG-CD	2.34	105.51	110.30	5	4
2	C	210	HYP	OD1-CG-CD	2.29	105.62	110.30	5	2
2	C	204	HYP	CG-CB-CA	2.29	106.39	103.75	3	1
2	D	322	HYP	OD1-CG-CD	2.27	105.67	110.30	4	3
2	B	119	HYP	CG-CB-CA	2.17	106.26	103.75	3	1
2	D	316	HYP	OD1-CG-CD	2.11	106.00	110.30	5	2
2	C	210	HYP	O-C-CA	2.07	119.45	124.77	8	2
2	B	104	HYP	CG-CB-CA	2.04	106.11	103.75	3	1
2	D	319	HYP	OD1-CG-CD	2.03	106.15	110.30	7	2
2	C	207	HYP	OD1-CG-CD	2.02	106.17	110.30	9	1
2	B	107	HYP	CG-CB-CA	2.01	106.08	103.75	3	1
2	D	304	HYP	OD1-CG-CD	2.00	106.21	110.30	9	1

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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

The completeness of assignment taking into account all chemical shift lists is 70% for the well-defined parts and 70% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_1*

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	2148
Number of shifts mapped to atoms	2148
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	4

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	191	-0.34 ± 0.14	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	178	-0.08 ± 0.13	None needed (< 0.5 ppm)
$^{13}\text{C}'$	157	0.00 ± 0.12	None needed (< 0.5 ppm)
^{15}N	181	0.91 ± 0.19	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 70%, i.e. 2058 atoms were assigned a chemical shift out of a possible 2951. 0 out of 34 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	869/1132 (77%)	363/469 (77%)	333/454 (73%)	173/209 (83%)
Sidechain	1130/1678 (67%)	771/1088 (71%)	352/528 (67%)	7/62 (11%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	59/141 (42%)	32/69 (46%)	26/63 (41%)	1/9 (11%)
Overall	2058/2951 (70%)	1166/1626 (72%)	711/1045 (68%)	181/280 (65%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 70%, i.e. 2148 atoms were assigned a chemical shift out of a possible 3074. 0 out of 35 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	909/1183 (77%)	380/490 (78%)	348/474 (73%)	181/219 (83%)
Sidechain	1174/1741 (67%)	799/1128 (71%)	367/546 (67%)	8/67 (12%)
Aromatic	65/150 (43%)	36/73 (49%)	28/68 (41%)	1/9 (11%)
Overall	2148/3074 (70%)	1215/1691 (72%)	743/1088 (68%)	190/295 (64%)

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	75	GLN	HB2	-0.19	0.80 – 3.29	-9.0
1	A	165	SER	HB2	1.99	2.61 – 5.13	-7.5
1	A	59	LYS	HG3	-0.24	0.04 – 2.67	-6.1
1	A	116	GLU	HG2	1.10	1.24 – 3.30	-5.7

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:

