



Full wwPDB NMR Structure Validation Report ⓘ

Apr 27, 2016 – 12:34 AM BST

PDB ID : 2L39
Title : Mouse prion protein fragment 121-231 AT 37 C
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Deposited on : 2010-09-10

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
<http://wwpdb.org/validation/2016/NMRValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

Cyrange : Kirchner and Güntert (2011)
NmrClust : Kelley et al. (1996)
MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : rb-20027457
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027457

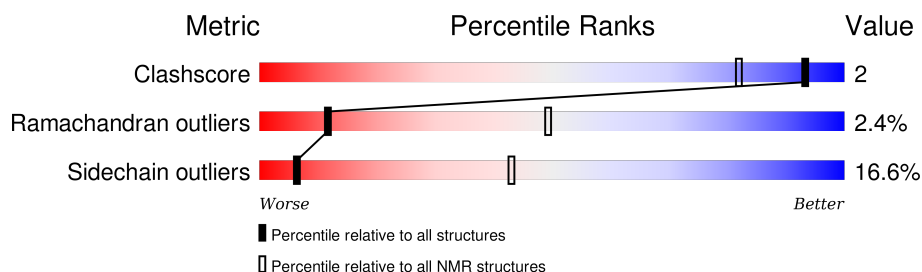
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR


The overall completeness of chemical shifts assignment is 82%.

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	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

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	114	 73% 16% 11%

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: *fewest violations*.

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:125-A:225 (101)	0.30	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 4 clusters and 2 single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1809 atoms, of which 872 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Major prion protein.

Mol	Chain	Residues	Atoms							Trace
1	A	114	Total	C	H	N	O	S		0
			1809	580	872	165	183	9		

There are 2 discrepancies between the modelled and reference sequences:

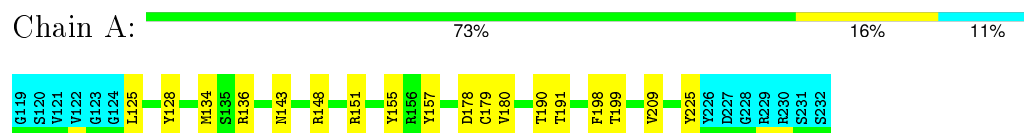
Chain	Residue	Modelled	Actual	Comment	Reference
A	119	GLY	-	SEE REMARK 999	UNP Q4FJQ7
A	120	SER	-	SEE REMARK 999	UNP Q4FJQ7

4 Residue-property plots [i](#)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA and DNA 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: Major prion protein

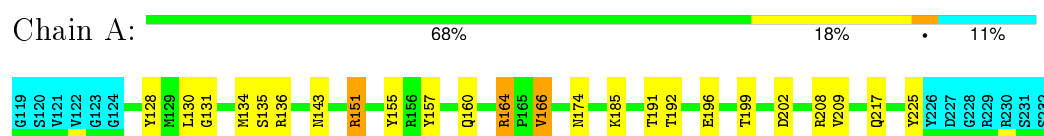


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

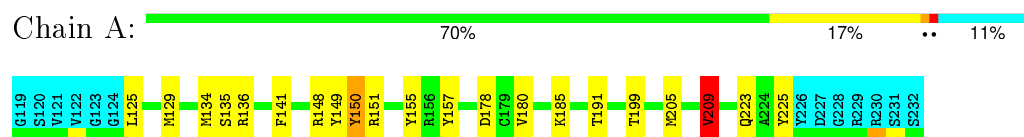
4.2.1 Score per residue for model 1

- Molecule 1: Major prion protein



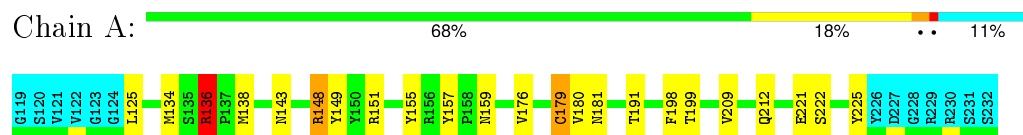
4.2.2 Score per residue for model 2

- Molecule 1: Major prion protein



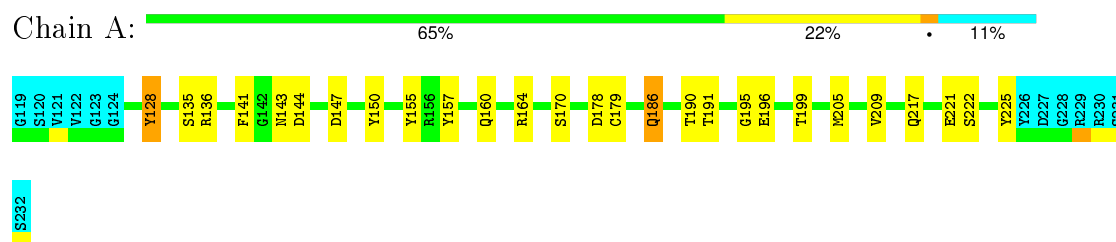
4.2.3 Score per residue for model 3 (medoid)

- Molecule 1: Major prion protein



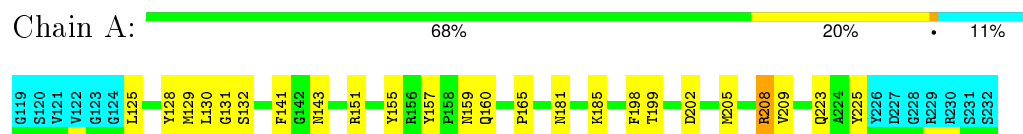
4.2.4 Score per residue for model 4

- Molecule 1: Major prion protein



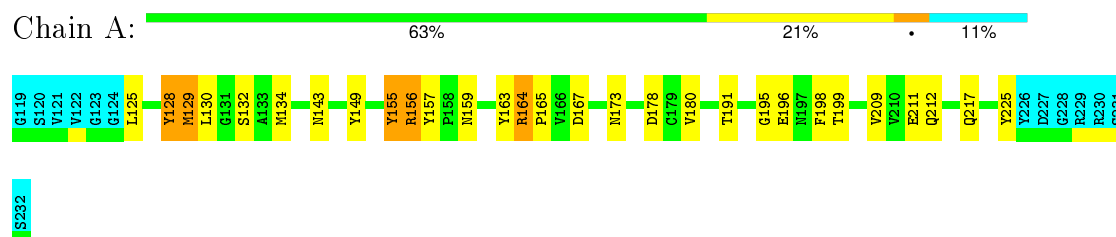
4.2.5 Score per residue for model 5

- Molecule 1: Major prion protein



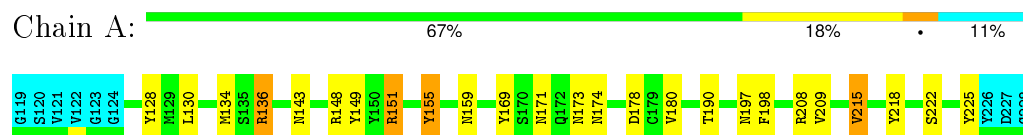
4.2.6 Score per residue for model 6

- Molecule 1: Major prion protein



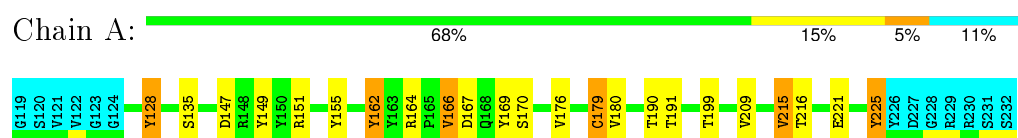
4.2.7 Score per residue for model 7

- Molecule 1: Major prion protein



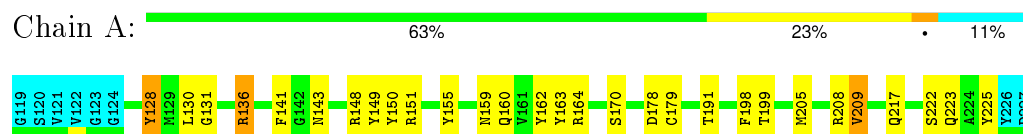
4.2.8 Score per residue for model 8

- Molecule 1: Major prion protein



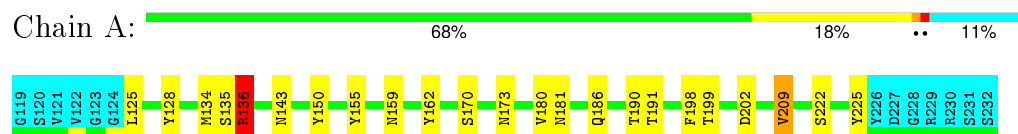
4.2.9 Score per residue for model 9

- Molecule 1: Major prion protein



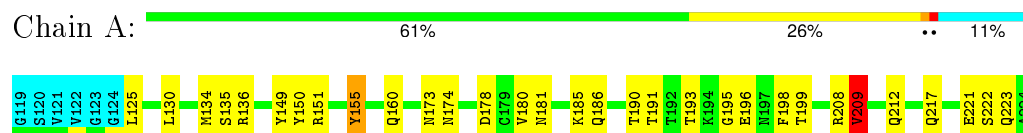
4.2.10 Score per residue for model 10

- Molecule 1: Major prion protein



4.2.11 Score per residue for model 11

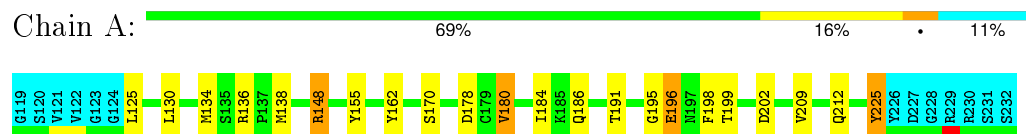
- Molecule 1: Major prion protein





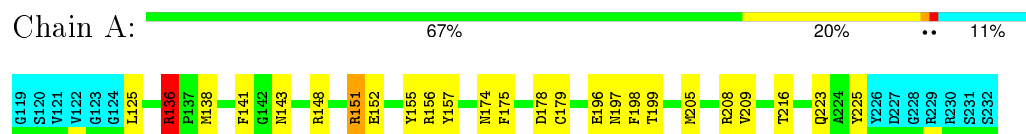
4.2.12 Score per residue for model 12

- Molecule 1: Major prion protein



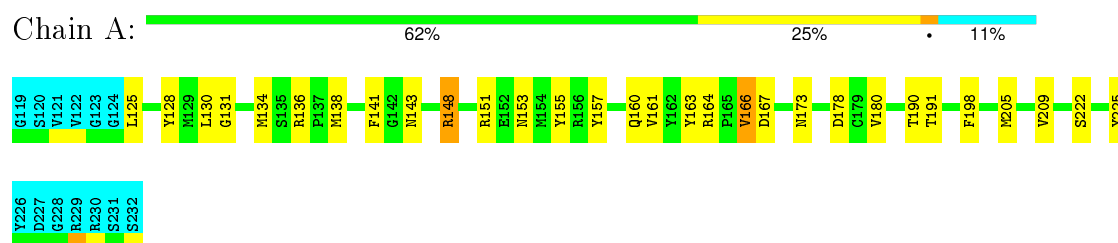
4.2.13 Score per residue for model 13

- Molecule 1: Major prion protein



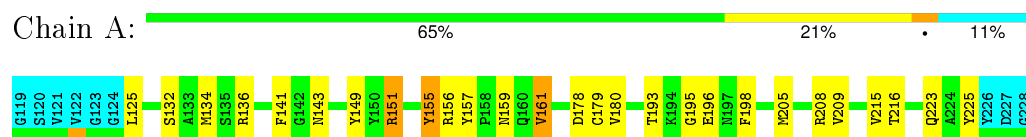
4.2.14 Score per residue for model 14

- Molecule 1: Major prion protein



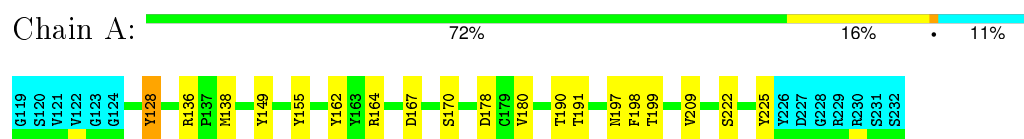
4.2.15 Score per residue for model 15

- Molecule 1: Major prion protein



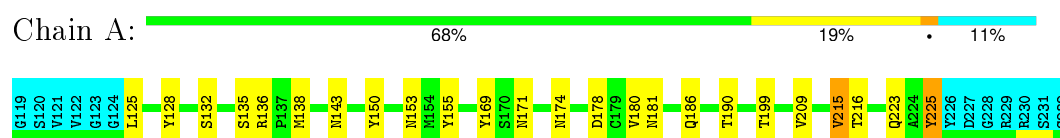
4.2.16 Score per residue for model 16

- Molecule 1: Major prion protein



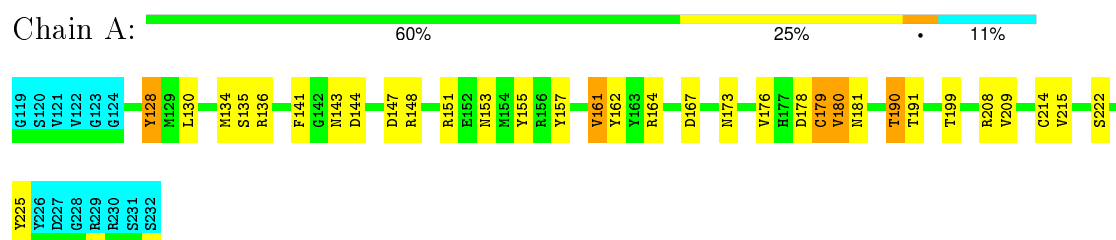
4.2.17 Score per residue for model 17

- Molecule 1: Major prion protein



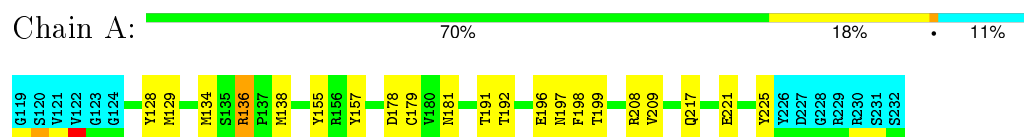
4.2.18 Score per residue for model 18

- Molecule 1: Major prion protein



4.2.19 Score per residue for model 19

- Molecule 1: Major prion protein



4.2.20 Score per residue for model 20

- Molecule 1: Major prion protein



G119	S120	V121	V122	G123	G124	L125	Y128	S132	R136	M143	R148	Y149	Y150	R151	Y155	R156	Y157	Y166	D167	Q168	Y169	M174	F175	V176	H177	D178	C179	T199	V209	Q217	Y225	Y226	D227	G228	R229	R230	S231	G232
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5 Refinement protocol and experimental data overview

The models were refined using the following method: *TORSION ANGLE DYNAMICS*,.

Of the 120 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
DYANA	structure solution	1.0.3
OPALP	refinement	1.2

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)	BMRB entry 17174
Number of chemical shift lists	1
Total number of shifts	1294
Number of shifts mapped to atoms	1294
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	82%

No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality i

6.1 Standard geometry i

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	0.68±0.01	0±0/868 (0.0±0.0%)	1.13±0.03	1±1/1175 (0.1±0.1%)
All	All	0.68	0/17360 (0.0%)	1.13	29/23500 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	3.9±1.7
All	All	0	77

There are no bond-length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	209	VAL	CA-CB-CG2	12.57	129.76	110.90	2	2
1	A	136	ARG	NE-CZ-NH2	-7.96	116.32	120.30	10	5
1	A	215	VAL	CA-CB-CG1	7.02	121.44	110.90	17	3
1	A	225	TYR	CB-CG-CD2	-7.02	116.79	121.00	8	1
1	A	156	ARG	NE-CZ-NH2	-6.98	116.81	120.30	6	2
1	A	161	VAL	CA-CB-CG1	5.95	119.83	110.90	15	2
1	A	166	VAL	CG1-CB-CG2	5.90	120.34	110.90	14	3
1	A	151	ARG	NE-CZ-NH2	-5.86	117.37	120.30	13	2
1	A	150	TYR	CB-CG-CD1	-5.59	117.64	121.00	2	2
1	A	208	ARG	NE-CZ-NH2	-5.52	117.54	120.30	11	2
1	A	148	ARG	NE-CZ-NH2	-5.51	117.55	120.30	12	1
1	A	151	ARG	NE-CZ-NH1	5.47	123.04	120.30	1	1
1	A	147	ASP	CB-CG-OD2	-5.24	113.58	118.30	18	1
1	A	162	TYR	CB-CG-CD1	-5.13	117.92	121.00	8	1
1	A	166	VAL	CA-CB-CG1	5.12	118.58	110.90	8	1

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	A	157	TYR	Sidechain	12
1	A	128	TYR	Sidechain	12
1	A	149	TYR	Sidechain	9
1	A	136	ARG	Sidechain	7
1	A	155	TYR	Sidechain	5
1	A	148	ARG	Sidechain	4
1	A	208	ARG	Sidechain	4
1	A	162	TYR	Sidechain	4
1	A	169	TYR	Sidechain	4
1	A	151	ARG	Sidechain	3
1	A	164	ARG	Sidechain	3
1	A	163	TYR	Sidechain	2
1	A	225	TYR	Sidechain	2
1	A	156	ARG	Sidechain	2
1	A	150	TYR	Sidechain	2
1	A	196	GLU	Peptide	1
1	A	218	TYR	Sidechain	1

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	847	788	788	3±2
All	All	16940	15760	15760	56

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:180:VAL:HG13	1:A:184:ILE:CD1	0.67	2.19	12	1
1:A:125:LEU:HD11	1:A:186:GLN:CD	0.60	2.17	10	1
1:A:180:VAL:HG13	1:A:184:ILE:HD12	0.56	1.77	12	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:130:LEU:HD13	1:A:162:TYR:CZ	0.56	2.35	18	1
1:A:176:VAL:HA	1:A:179:CYS:SG	0.55	2.42	3	4
1:A:150:TYR:CE2	1:A:209:VAL:HG13	0.52	2.40	2	1
1:A:130:LEU:HD11	1:A:160:GLN:HE21	0.51	1.66	14	1
1:A:130:LEU:HD22	1:A:131:GLY:H	0.50	1.65	14	2
1:A:150:TYR:CE1	1:A:209:VAL:HG13	0.50	2.41	11	1
1:A:161:VAL:HG21	1:A:214:CYS:SG	0.49	2.46	18	1
1:A:171:ASN:ND2	1:A:174:ASN:H	0.49	2.05	17	1
1:A:150:TYR:OH	1:A:209:VAL:HG23	0.49	2.07	10	1
1:A:186:GLN:O	1:A:190:THR:HG22	0.49	2.07	4	1
1:A:141:PHE:CE2	1:A:205:MET:SD	0.48	3.07	9	7
1:A:128:TYR:CE2	1:A:164:ARG:HG2	0.47	2.44	4	4
1:A:130:LEU:HD13	1:A:131:GLY:N	0.47	2.25	1	2
1:A:180:VAL:CG1	1:A:184:ILE:HD12	0.46	2.40	12	1
1:A:125:LEU:HD21	1:A:186:GLN:CD	0.46	2.31	17	3
1:A:130:LEU:HD11	1:A:160:GLN:HB2	0.45	1.88	11	1
1:A:171:ASN:HD21	1:A:174:ASN:H	0.45	1.54	17	1
1:A:130:LEU:HD21	1:A:160:GLN:HB2	0.45	1.88	9	1
1:A:150:TYR:CE1	1:A:209:VAL:HG23	0.45	2.47	9	1
1:A:129:MET:CE	1:A:163:TYR:CE2	0.44	3.00	6	1
1:A:171:ASN:HD21	1:A:174:ASN:HB2	0.44	1.73	7	1
1:A:179:CYS:SG	1:A:180:VAL:N	0.44	2.90	18	1
1:A:130:LEU:HD22	1:A:161:VAL:O	0.44	2.13	14	1
1:A:130:LEU:HD13	1:A:162:TYR:CE1	0.44	2.48	12	1
1:A:196:GLU:CD	1:A:198:PHE:CD2	0.43	2.92	12	1
1:A:130:LEU:HD11	1:A:160:GLN:NE2	0.42	2.30	1	2
1:A:128:TYR:CZ	1:A:164:ARG:HD3	0.42	2.49	9	4
1:A:141:PHE:CE2	1:A:208:ARG:HG3	0.42	2.50	18	1
1:A:129:MET:SD	1:A:165:PRO:HD3	0.41	2.55	6	1
1:A:191:THR:HG22	1:A:196:GLU:CB	0.41	2.45	4	1
1:A:175:PHE:CE1	1:A:179:CYS:SG	0.41	3.13	13	1
1:A:196:GLU:HG3	1:A:198:PHE:CE2	0.40	2.51	6	1
1:A:150:TYR:HE2	1:A:209:VAL:HG13	0.40	1.74	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	101/114 (89%)	88±3 (87±3%)	11±3 (11±3%)	2±1 (2±1%)	12	49
All	All	2020/2280 (89%)	1758 (87%)	213 (11%)	49 (2%)	12	49

All 12 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	180	VAL	13
1	A	198	PHE	11
1	A	125	LEU	8
1	A	195	GLY	5
1	A	190	THR	3
1	A	170	SER	2
1	A	197	ASN	2
1	A	167	ASP	1
1	A	128	TYR	1
1	A	143	ASN	1
1	A	165	PRO	1
1	A	169	TYR	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	94/103 (91%)	78±3 (83±3%)	16±3 (17±3%)	6	43
All	All	1880/2060 (91%)	1567 (83%)	313 (17%)	6	43

All 47 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	209	VAL	20
1	A	225	TYR	20
1	A	155	TYR	20
1	A	199	THR	17
1	A	178	ASP	15

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Mol	Chain	Res	Type	Models (Total)
1	A	136	ARG	14
1	A	191	THR	13
1	A	143	ASN	13
1	A	134	MET	12
1	A	151	ARG	11
1	A	222	SER	9
1	A	148	ARG	8
1	A	135	SER	8
1	A	181	ASN	7
1	A	159	ASN	7
1	A	179	CYS	7
1	A	217	GLN	7
1	A	138	MET	7
1	A	223	GLN	7
1	A	190	THR	6
1	A	173	ASN	6
1	A	167	ASP	5
1	A	132	SER	5
1	A	221	GLU	5
1	A	196	GLU	5
1	A	215	VAL	5
1	A	202	ASP	4
1	A	185	LYS	4
1	A	166	VAL	4
1	A	174	ASN	4
1	A	212	GLN	4
1	A	216	THR	4
1	A	170	SER	4
1	A	208	ARG	3
1	A	129	MET	3
1	A	153	ASN	3
1	A	193	THR	2
1	A	147	ASP	2
1	A	192	THR	2
1	A	144	ASP	2
1	A	130	LEU	2
1	A	197	ASN	2
1	A	160	GLN	1
1	A	186	GLN	1
1	A	211	GLU	1
1	A	161	VAL	1
1	A	152	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 carbohydrates 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 82% for the well-defined parts and 81% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 17174

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	1294
Number of shifts mapped to atoms	1294
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	114	-0.41 ± 0.17	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	105	0.24 ± 0.13	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	—
^{15}N	110	0.13 ± 0.17	None needed (< 0.5 ppm)

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 82%, i.e. 1059 atoms were assigned a chemical shift out of a possible 1295. 11 out of 11 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	392/499 (79%)	193/199 (97%)	101/202 (50%)	98/98 (100%)
Sidechain	563/656 (86%)	353/387 (91%)	188/231 (81%)	22/38 (58%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	104/140 (74%)	66/73 (90%)	37/63 (59%)	1/4 (25%)
Overall	1059/1295 (82%)	612/659 (93%)	326/496 (66%)	121/140 (86%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 81%, i.e. 1164 atoms were assigned a chemical shift out of a possible 1432. 13 out of 13 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	442/564 (78%)	218/225 (97%)	114/228 (50%)	110/111 (99%)
Sidechain	612/720 (85%)	383/425 (90%)	205/251 (82%)	24/44 (55%)
Aromatic	110/148 (74%)	70/77 (91%)	39/67 (58%)	1/4 (25%)
Overall	1164/1432 (81%)	671/727 (92%)	358/546 (66%)	135/159 (85%)

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.

Random coil index (RCI) for chain A:

