



Full wwPDB NMR Structure Validation Report ⓘ

Feb 12, 2017 – 09:01 pm GMT

PDB ID : 2AGM
Title : Solution structure of the R-module from AlgE4
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Deposited on : 2005-07-27

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
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : trunk28760
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

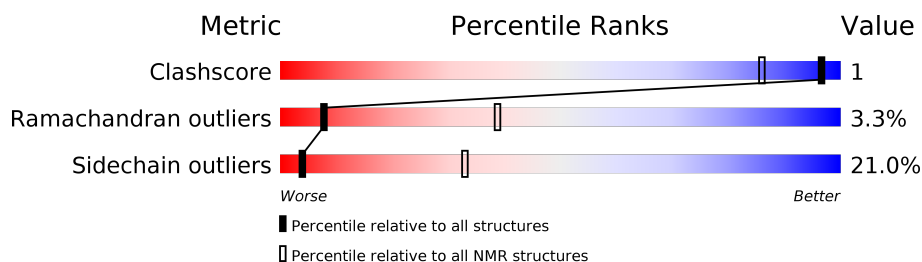
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 81%.

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	125131	11601
Ramachandran outliers	121729	10391
Sidechain outliers	121581	10367

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	167	

2 Ensemble composition and analysis

This entry contains 20 models. Model 2 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:144 (141)	0.87	2

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 3 clusters and 3 single-model clusters were found.

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

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 2296 atoms, of which 1102 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Poly(beta-D-mannuronate) C5 epimerase 4.

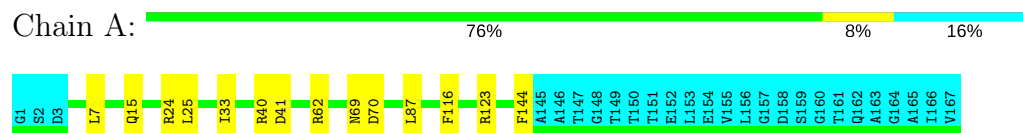
Mol	Chain	Residues	Atoms					Trace
1	A	167	Total	C	H	N	O	0
			2296	723	1102	204	267	

4 Residue-property plots

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: Poly(beta-D-mannuronate) C5 epimerase 4

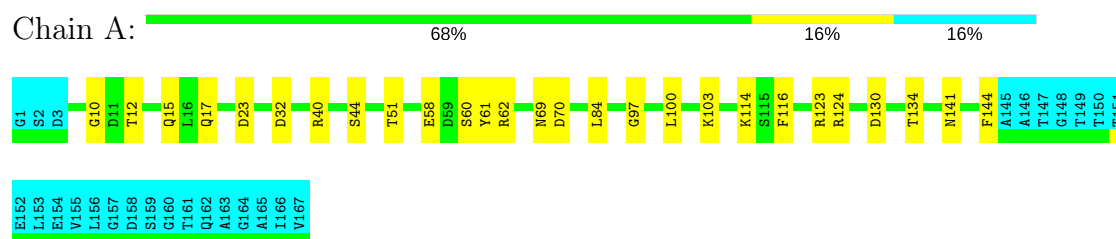


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

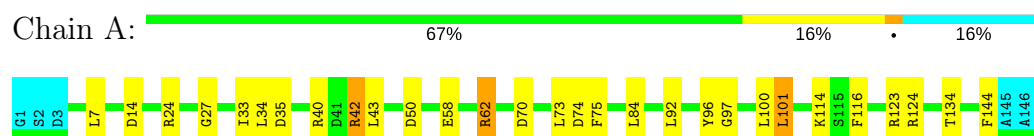
4.2.1 Score per residue for model 1

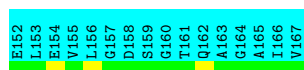
- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



4.2.2 Score per residue for model 2 (medoid)

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4

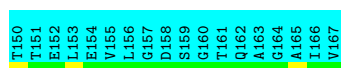
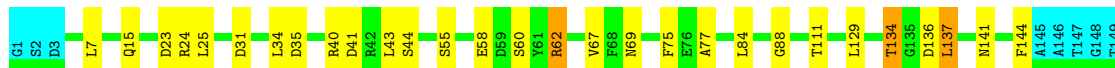




4.2.3 Score per residue for model 3

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4

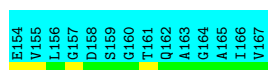
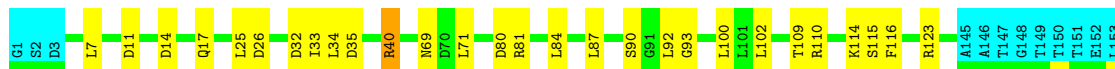
Chain A: 67% 16% 16%



4.2.4 Score per residue for model 4

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4

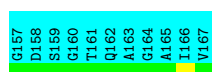
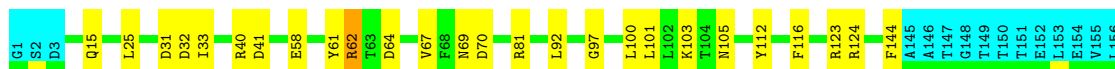
Chain A: 68% 16% 16%



4.2.5 Score per residue for model 5

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4

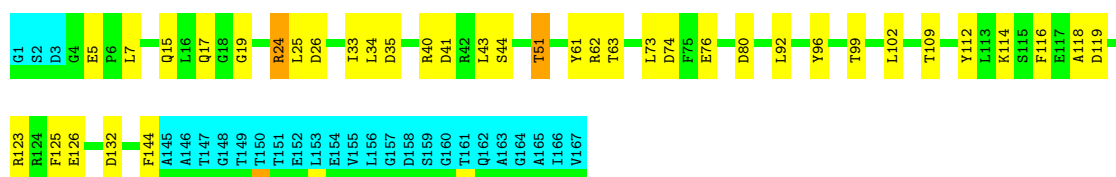
Chain A: 69% 15% 16%



4.2.6 Score per residue for model 6

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4

Chain A: 62% 22% 16%



4.2.7 Score per residue for model 7

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4

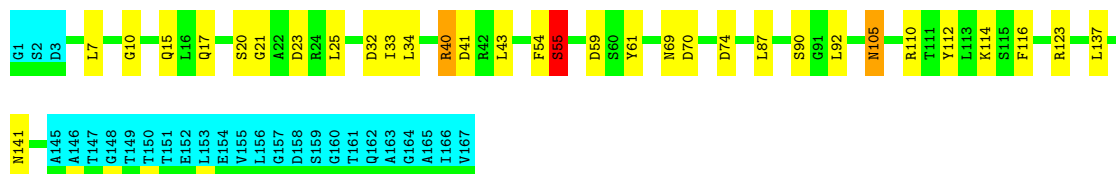
Chain A: 69% 14% 16%



4.2.8 Score per residue for model 8

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4

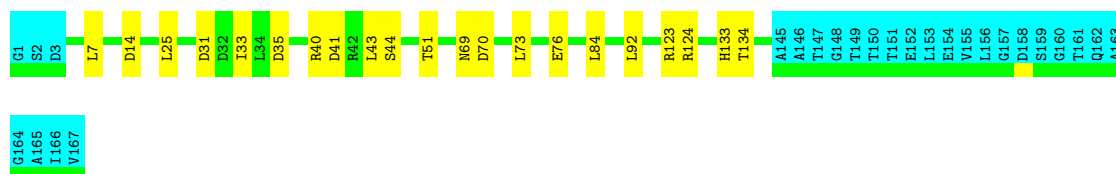
Chain A: 65% 17% 16%



4.2.9 Score per residue for model 9

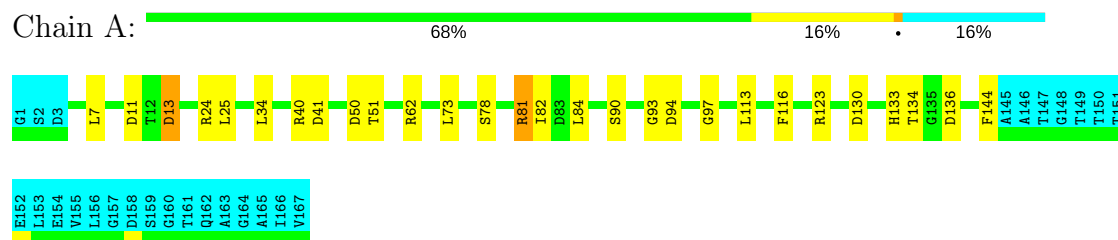
- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4

Chain A: 72% 13% 16%



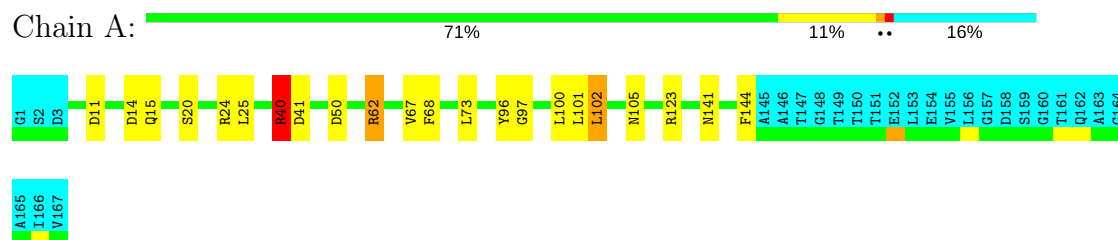
4.2.10 Score per residue for model 10

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



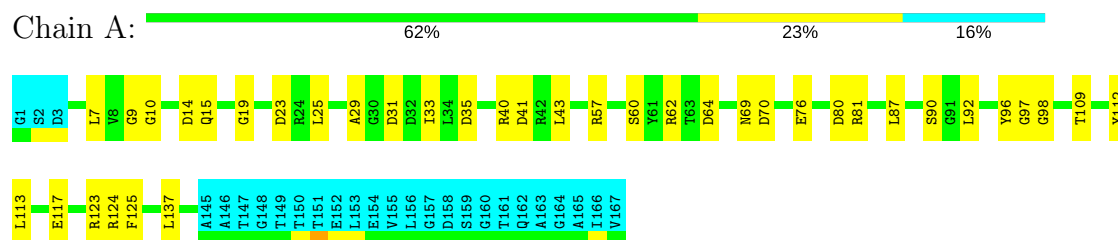
4.2.11 Score per residue for model 11

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



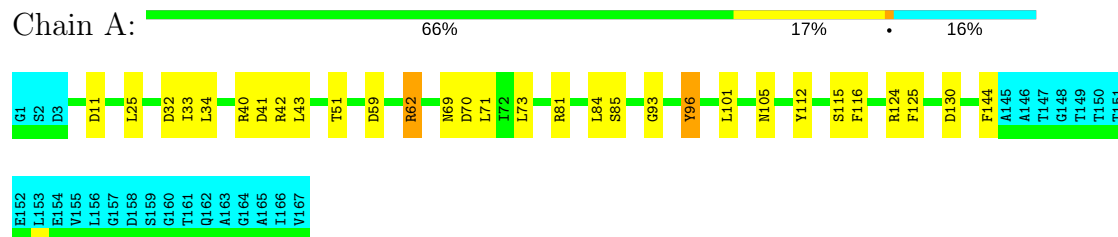
4.2.12 Score per residue for model 12

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



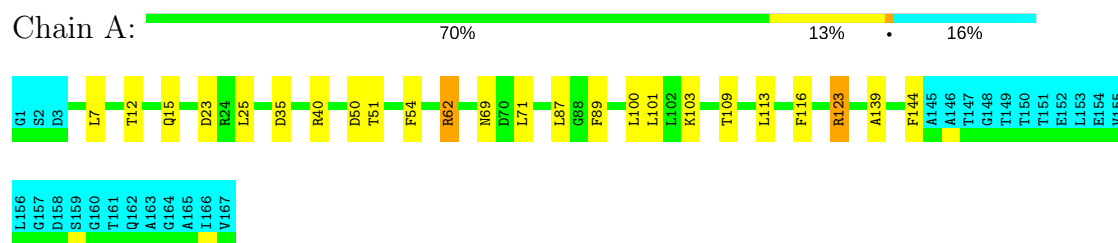
4.2.13 Score per residue for model 13

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



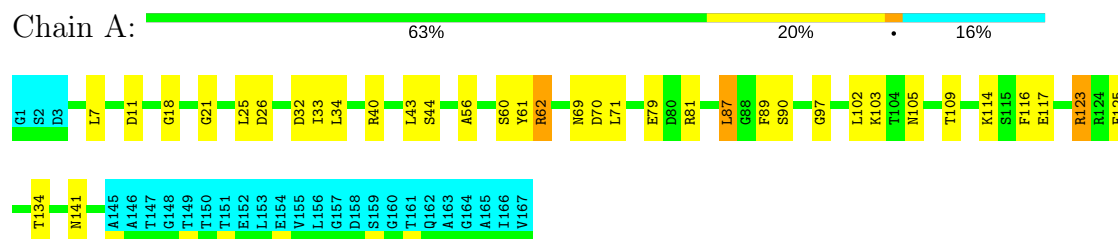
4.2.14 Score per residue for model 14

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



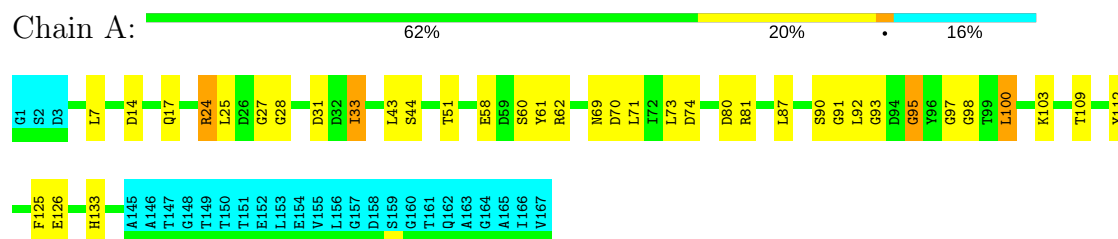
4.2.15 Score per residue for model 15

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



4.2.16 Score per residue for model 16

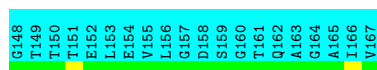
- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



4.2.17 Score per residue for model 17

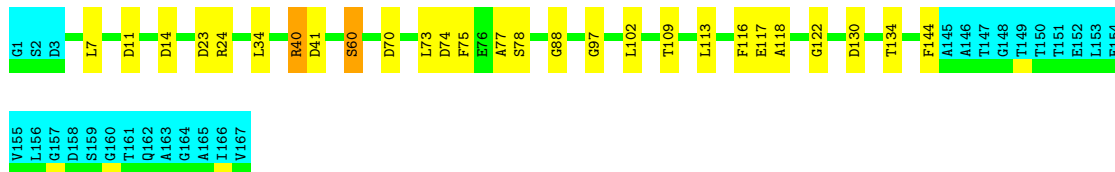
- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4





4.2.18 Score per residue for model 18

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



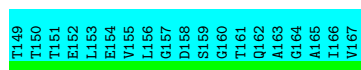
4.2.19 Score per residue for model 19

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



4.2.20 Score per residue for model 20

- Molecule 1: Poly(beta-D-mannuronate) C5 epimerase 4



5 Refinement protocol and experimental data overview

The models were refined using the following method: *DISTANCE GEOMETRY, torsion angle dynamics, energy-minimisation*.

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	structure solution	2.1.5
OPALp	structure solution	1.4
OPALp	refinement	1.4

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 6390
Number of chemical shift lists	1
Total number of shifts	1661
Number of shifts mapped to atoms	1661
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	81%

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	0.61±0.01	0±0/1039 (0.0±0.0%)	1.19±0.03	2±1/1402 (0.1±0.1%)
All	All	0.61	0/20780 (0.0%)	1.19	40/28040 (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	1.1±0.8
All	All	0	22

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	62	ARG	NE-CZ-NH1	7.15	123.87	120.30	15	2
1	A	59	ASP	C-N-CA	6.71	138.48	121.70	19	1
1	A	62	ARG	NE-CZ-NH2	-6.66	116.97	120.30	14	2
1	A	40	ARG	NE-CZ-NH2	-6.47	117.07	120.30	11	3
1	A	11	ASP	CB-CG-OD1	6.17	123.85	118.30	4	1
1	A	31	ASP	N-CA-CB	-6.08	99.66	110.60	3	1
1	A	62	ARG	CD-NE-CZ	6.07	132.10	123.60	15	1
1	A	40	ARG	CB-CG-CD	6.05	127.32	111.60	11	1
1	A	89	PHE	C-N-CA	5.93	136.52	121.70	7	1
1	A	23	ASP	CB-CA-C	5.90	122.19	110.40	8	1
1	A	31	ASP	CB-CA-C	5.79	121.98	110.40	9	1
1	A	96	TYR	C-N-CA	5.78	134.44	122.30	12	2
1	A	134	THR	CA-CB-CG2	5.71	120.39	112.40	3	2
1	A	23	ASP	CB-CG-OD2	5.70	123.43	118.30	8	1
1	A	35	ASP	CB-CG-OD1	5.69	123.42	118.30	3	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	124	ARG	NE-CZ-NH2	-5.69	117.46	120.30	1	1
1	A	124	ARG	NE-CZ-NH1	5.61	123.11	120.30	5	1
1	A	55	SER	CB-CA-C	5.51	120.57	110.10	8	1
1	A	116	PHE	N-CA-CB	-5.40	100.87	110.60	20	1
1	A	24	ARG	NE-CZ-NH2	-5.37	117.62	120.30	17	1
1	A	110	ARG	NE-CZ-NH2	-5.29	117.66	120.30	4	1
1	A	59	ASP	CB-CG-OD2	-5.25	113.58	118.30	13	1
1	A	125	PHE	N-CA-CB	-5.22	101.21	110.60	6	1
1	A	33	ILE	CA-CB-CG1	5.22	120.92	111.00	20	1
1	A	118	ALA	CB-CA-C	5.12	117.78	110.10	19	1
1	A	123	ARG	NE-CZ-NH2	-5.07	117.76	120.30	12	1
1	A	31	ASP	CB-CG-OD1	5.05	122.85	118.30	16	1
1	A	51	THR	CA-CB-CG2	5.04	119.46	112.40	6	1
1	A	112	TYR	CB-CG-CD2	-5.04	117.98	121.00	13	1
1	A	81	ARG	NE-CZ-NH1	5.03	122.82	120.30	13	1
1	A	73	LEU	C-N-CA	5.03	134.27	121.70	6	1
1	A	27	GLY	C-N-CA	5.03	132.86	122.30	19	1
1	A	117	GLU	C-N-CA	5.01	134.23	121.70	20	1
1	A	13	ASP	CB-CG-OD1	5.00	122.80	118.30	10	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	62	ARG	Sidechain	4
1	A	40	ARG	Sidechain	4
1	A	124	ARG	Sidechain	3
1	A	21	GLY	Mainchain	2
1	A	24	ARG	Sidechain	2
1	A	112	TYR	Sidechain	1
1	A	81	ARG	Sidechain	1
1	A	123	ARG	Sidechain	1
1	A	125	PHE	Sidechain	1
1	A	42	ARG	Sidechain	1
1	A	61	TYR	Sidechain	1
1	A	57	ARG	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	1026	943	943	2±1
All	All	20520	18860	18860	35

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 1.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:55:SER:HA	1:A:87:LEU:HD13	0.59	1.75	8	1
1:A:40:ARG:HG3	1:A:68:PHE:HA	0.51	1.83	11	1
1:A:77:ALA:HB3	1:A:134:THR:O	0.49	2.07	3	1
1:A:40:ARG:CG	1:A:68:PHE:HA	0.49	2.37	11	1
1:A:15:GLN:HG2	1:A:33:ILE:CG2	0.49	2.38	8	3
1:A:27:GLY:N	1:A:34:LEU:HD21	0.48	2.24	2	1
1:A:60:SER:OG	1:A:87:LEU:HD22	0.47	2.09	12	1
1:A:62:ARG:HB3	1:A:67:VAL:HG13	0.47	1.85	17	1
1:A:56:ALA:CB	1:A:87:LEU:HD13	0.47	2.40	15	1
1:A:81:ARG:C	1:A:82:ILE:HD12	0.46	2.31	10	1
1:A:75:PHE:CG	1:A:129:LEU:HD11	0.46	2.45	3	1
1:A:55:SER:HB2	1:A:87:LEU:HD11	0.46	1.86	17	1
1:A:95:GLY:HA3	1:A:100:LEU:HA	0.45	1.87	16	1
1:A:43:LEU:C	1:A:43:LEU:HD12	0.44	2.32	16	1
1:A:33:ILE:HG22	1:A:51:THR:HB	0.44	1.89	13	2
1:A:75:PHE:CE2	1:A:77:ALA:HB2	0.44	2.47	18	1
1:A:54:PHE:O	1:A:86:ALA:HB3	0.43	2.14	17	1
1:A:75:PHE:CD2	1:A:129:LEU:HD11	0.43	2.48	20	1
1:A:33:ILE:C	1:A:33:ILE:HD12	0.43	2.33	19	1
1:A:55:SER:CA	1:A:87:LEU:HD13	0.42	2.43	8	1
1:A:15:GLN:HG2	1:A:33:ILE:HD11	0.41	1.91	19	1
1:A:101:LEU:HD23	1:A:102:LEU:N	0.41	2.30	11	1
1:A:75:PHE:CE2	1:A:129:LEU:HD11	0.41	2.50	20	1
1:A:105:ASN:ND2	1:A:110:ARG:H	0.41	2.14	8	1
1:A:63:THR:HG22	1:A:123:ARG:HG3	0.41	1.92	7	1
1:A:113:LEU:H	1:A:113:LEU:HD23	0.41	1.75	14	1
1:A:100:LEU:HD23	1:A:101:LEU:N	0.41	2.31	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:15:GLN:HG3	1:A:33:ILE:CG1	0.40	2.47	19	1
1:A:56:ALA:HA	1:A:86:ALA:HB1	0.40	1.92	17	1
1:A:137:LEU:HA	1:A:141:ASN:HD22	0.40	1.77	3	1
1:A:54:PHE:HB3	1:A:87:LEU:HD23	0.40	1.92	14	1
1:A:119:ASP:HB3	1:A:123:ARG:H	0.40	1.77	20	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	141/167 (84%)	115±3 (82±2%)	21±3 (15±2%)	5±2 (3±2%)	8	39
All	All	2820/3340 (84%)	2304 (82%)	422 (15%)	94 (3%)	8	39

All 37 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	116	PHE	13
1	A	97	GLY	10
1	A	123	ARG	8
1	A	93	GLY	5
1	A	74	ASP	4
1	A	64	ASP	4
1	A	60	SER	4
1	A	11	ASP	3
1	A	88	GLY	3
1	A	10	GLY	3
1	A	118	ALA	3
1	A	96	TYR	2
1	A	28	GLY	2
1	A	18	GLY	2
1	A	94	ASP	2
1	A	98	GLY	2
1	A	55	SER	2

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Mol	Chain	Res	Type	Models (Total)
1	A	19	GLY	2
1	A	57	ARG	2
1	A	89	PHE	1
1	A	122	GLY	1
1	A	139	ALA	1
1	A	95	GLY	1
1	A	90	SER	1
1	A	20	SER	1
1	A	56	ALA	1
1	A	9	GLY	1
1	A	92	LEU	1
1	A	36	GLY	1
1	A	133	HIS	1
1	A	29	ALA	1
1	A	134	THR	1
1	A	91	GLY	1
1	A	27	GLY	1
1	A	135	GLY	1
1	A	117	GLU	1
1	A	5	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	101/118 (86%)	80±4 (79±4%)	21±4 (21±4%)	4	33
All	All	2020/2360 (86%)	1595 (79%)	425 (21%)	4	33

All 85 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	40	ARG	15
1	A	7	LEU	14
1	A	144	PHE	14
1	A	62	ARG	13
1	A	70	ASP	13

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Mol	Chain	Res	Type	Models (Total)
1	A	25	LEU	13
1	A	41	ASP	12
1	A	69	ASN	12
1	A	43	LEU	9
1	A	109	THR	9
1	A	33	ILE	9
1	A	73	LEU	9
1	A	61	TYR	9
1	A	24	ARG	8
1	A	34	LEU	8
1	A	14	ASP	8
1	A	100	LEU	8
1	A	114	LYS	8
1	A	35	ASP	8
1	A	84	LEU	8
1	A	92	LEU	7
1	A	90	SER	7
1	A	17	GLN	7
1	A	15	GLN	7
1	A	58	GLU	7
1	A	44	SER	6
1	A	80	ASP	6
1	A	81	ARG	6
1	A	141	ASN	6
1	A	32	ASP	6
1	A	23	ASP	6
1	A	123	ARG	6
1	A	102	LEU	6
1	A	87	LEU	5
1	A	113	LEU	5
1	A	124	ARG	5
1	A	134	THR	5
1	A	103	LYS	5
1	A	105	ASN	5
1	A	71	LEU	5
1	A	125	PHE	5
1	A	50	ASP	5
1	A	51	THR	5
1	A	60	SER	4
1	A	11	ASP	4
1	A	112	TYR	4
1	A	130	ASP	4

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Mol	Chain	Res	Type	Models (Total)
1	A	101	LEU	4
1	A	126	GLU	3
1	A	136	ASP	3
1	A	76	GLU	3
1	A	137	LEU	3
1	A	26	ASP	3
1	A	133	HIS	3
1	A	115	SER	3
1	A	117	GLU	3
1	A	67	VAL	3
1	A	96	TYR	2
1	A	31	ASP	2
1	A	78	SER	2
1	A	20	SER	2
1	A	116	PHE	2
1	A	42	ARG	2
1	A	12	THR	2
1	A	57	ARG	2
1	A	74	ASP	2
1	A	82	ILE	2
1	A	132	ASP	1
1	A	99	THR	1
1	A	119	ASP	1
1	A	138	SER	1
1	A	94	ASP	1
1	A	110	ARG	1
1	A	127	VAL	1
1	A	55	SER	1
1	A	59	ASP	1
1	A	111	THR	1
1	A	79	GLU	1
1	A	13	ASP	1
1	A	75	PHE	1
1	A	54	PHE	1
1	A	89	PHE	1
1	A	129	LEU	1
1	A	85	SER	1
1	A	63	THR	1

6.3.3 RNA ⓘ

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 81% for the well-defined parts and 82% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 6390

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	1661
Number of shifts mapped to atoms	1661
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
$^{13}\text{C}_\alpha$	166	0.03 ± 0.14	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	137	0.31 ± 0.14	None needed (< 0.5 ppm)
$^{13}\text{C}'$	158	0.03 ± 0.13	None needed (< 0.5 ppm)
^{15}N	159	-1.11 ± 0.41	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 81%, i.e. 1234 atoms were assigned a chemical shift out of a possible 1520. 17 out of 22 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	681/703 (97%)	274/281 (98%)	273/282 (97%)	134/140 (96%)
Sidechain	532/714 (75%)	327/407 (80%)	199/273 (73%)	6/34 (18%)

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	Total	¹H	¹³C	¹⁵N
Aromatic	21/103 (20%)	21/56 (38%)	0/46 (0%)	0/1 (0%)
Overall	1234/1520 (81%)	622/744 (84%)	472/601 (79%)	140/175 (80%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 82%, i.e. 1447 atoms were assigned a chemical shift out of a possible 1755. 21 out of 26 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	808/833 (97%)	325/333 (98%)	324/334 (97%)	159/166 (96%)
Sidechain	618/819 (75%)	377/464 (81%)	234/320 (73%)	7/35 (20%)
Aromatic	21/103 (20%)	21/56 (38%)	0/46 (0%)	0/1 (0%)
Overall	1447/1755 (82%)	723/853 (85%)	558/700 (80%)	166/202 (82%)

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.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	57	ARG	NE	111.90	92.63 – 76.73	17.1
1	A	91	GLY	HA3	1.68	5.80 – 2.00	-5.8
1	A	33	ILE	CG1	18.10	36.54 – 18.94	-5.5
1	A	45	GLY	H	11.76	11.63 – 5.03	5.2

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.

Random coil index (RCI) for chain A:

