



Full wwPDB NMR Structure Validation Report ⓘ

Feb 12, 2017 – 10:15 pm GMT

PDB ID : 2JO0
Title : The solution structure of the monomeric species of the C terminal domain of the CA protein of HIV-1
Authors : Alcaraz, L.A.; del Alamo, M.; Barrera, F.N.; Mateu, M.G.; Neira, J.L.
Deposited on : 2007-02-17

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

<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

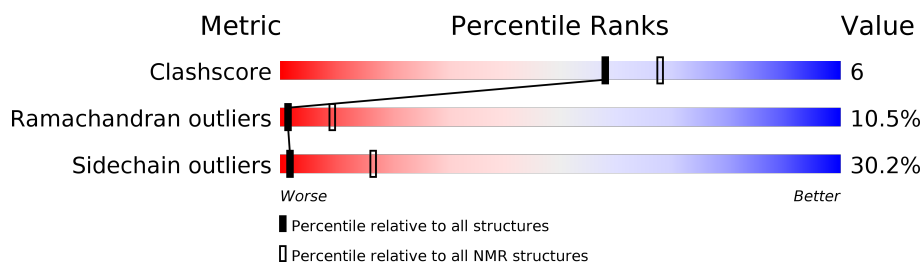
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 51%.

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	87	<div>51% 26% • 20%</div>

2 Ensemble composition and analysis

This entry contains 30 models. The atoms present in the NMR models are not consistent. Some calculations may have failed as a result. All residues are included in the validation scores. Model 14 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:6-A:75 (70)	0.84	14

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 4 single-model clusters were found.

Cluster number	Models
1	7, 12, 14, 16, 20, 21, 22, 27, 28
2	2, 3, 4, 8, 9, 25
3	1, 10, 13, 17, 24, 26
4	6, 15, 29
5	5, 11
Single-model clusters	18; 19; 23; 30

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1343 atoms, of which 677 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Gag-Pol polyprotein.

Mol	Chain	Residues	Atoms						Trace
1	A	87	Total	C	H	N	O	S	0
			1343	414	677	118	128	6	

There are 2 discrepancies between the modelled and reference sequences:

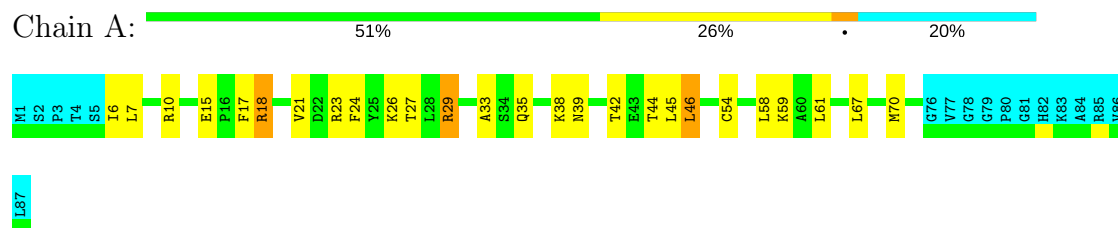
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP P35963
A	40	ALA	TRP	ENGINEERED	UNP P35963

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: Gag-Pol polyprotein

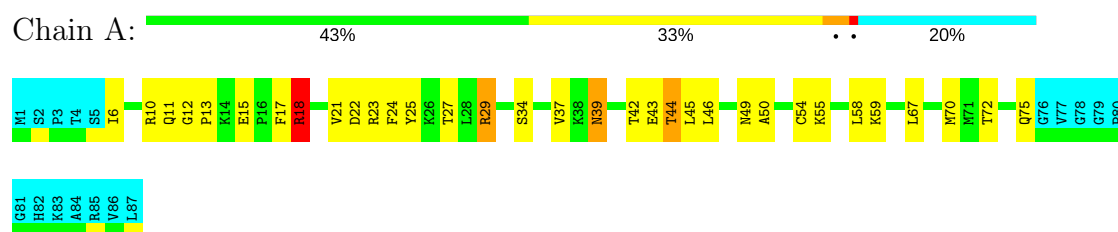


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

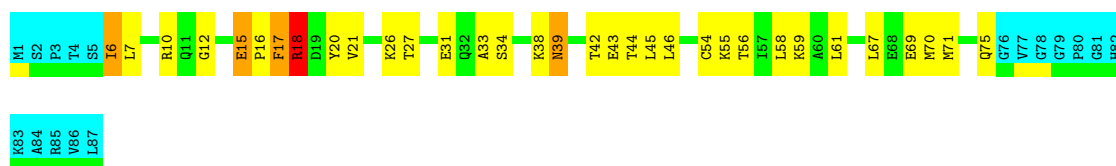
- Molecule 1: Gag-Pol polyprotein



4.2.2 Score per residue for model 2

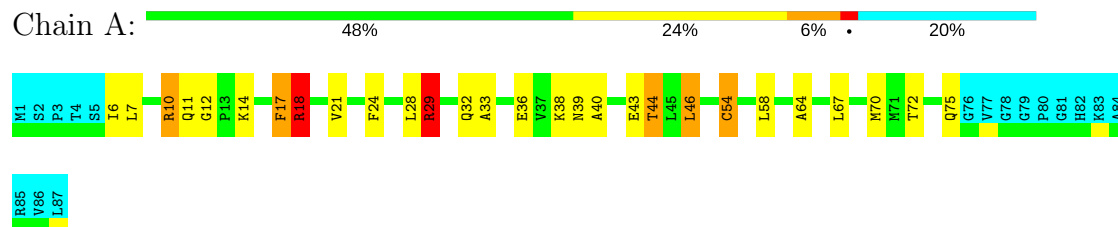
- Molecule 1: Gag-Pol polyprotein





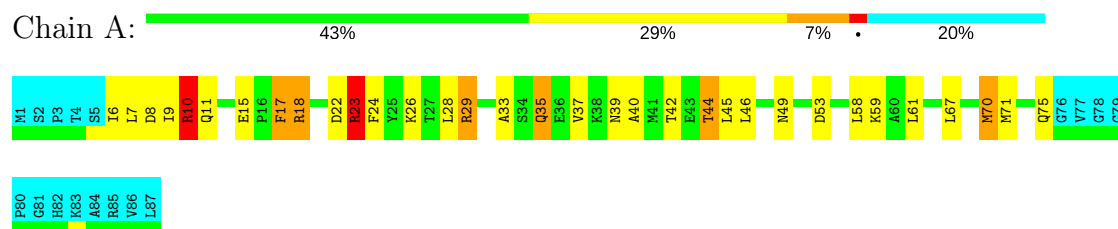
4.2.3 Score per residue for model 3

- Molecule 1: Gag-Pol polyprotein



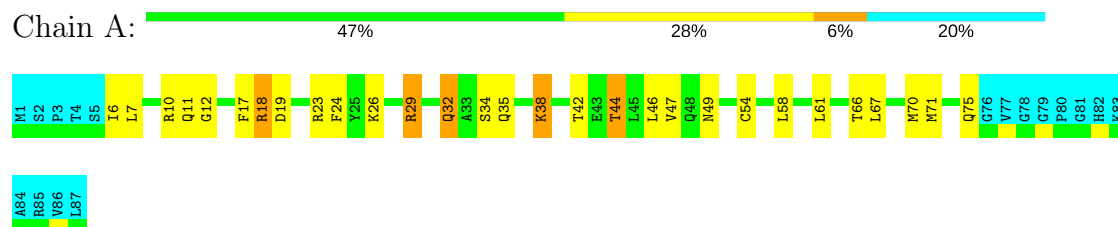
4.2.4 Score per residue for model 4

- Molecule 1: Gag-Pol polyprotein



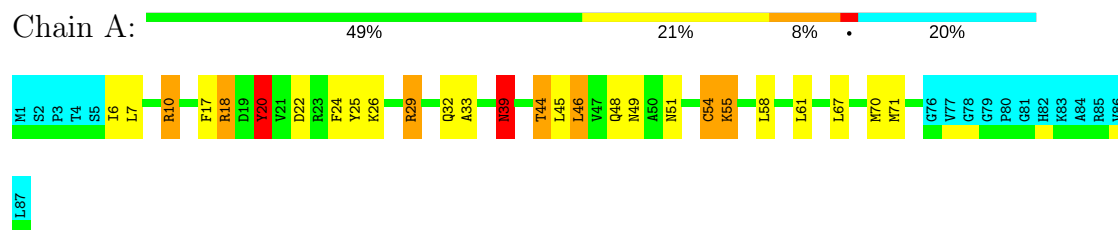
4.2.5 Score per residue for model 5

- Molecule 1: Gag-Pol polyprotein



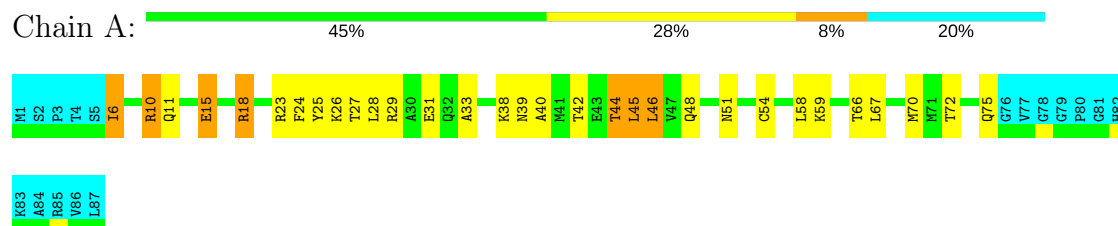
4.2.6 Score per residue for model 6

- Molecule 1: Gag-Pol polyprotein



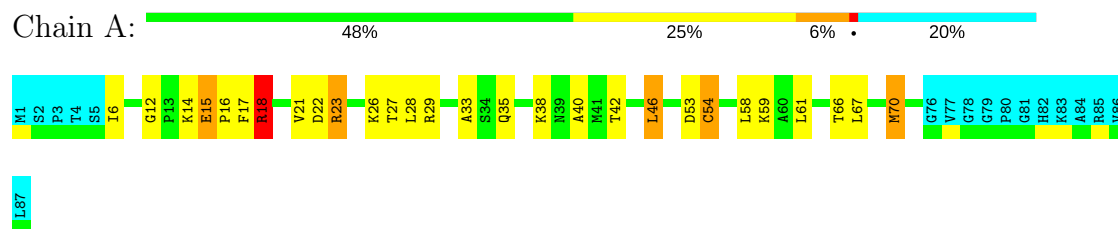
4.2.7 Score per residue for model 7

- Molecule 1: Gag-Pol polyprotein



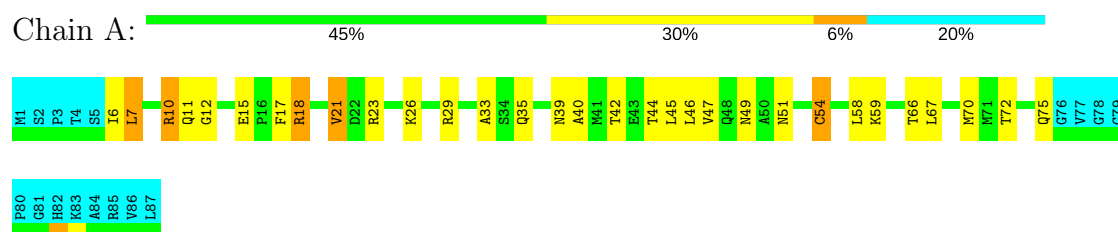
4.2.8 Score per residue for model 8

- Molecule 1: Gag-Pol polyprotein



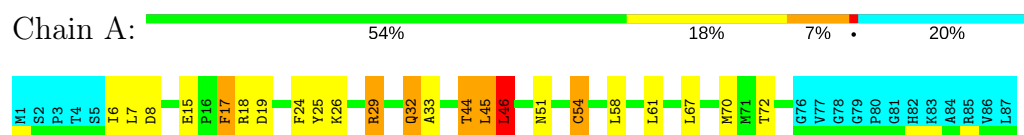
4.2.9 Score per residue for model 9

- Molecule 1: Gag-Pol polyprotein



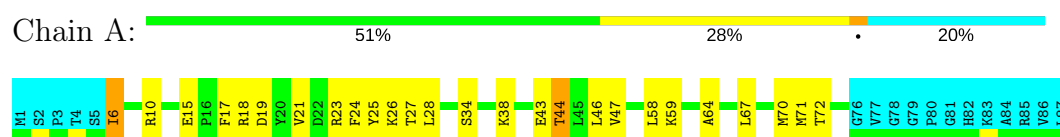
4.2.10 Score per residue for model 10

- Molecule 1: Gag-Pol polyprotein



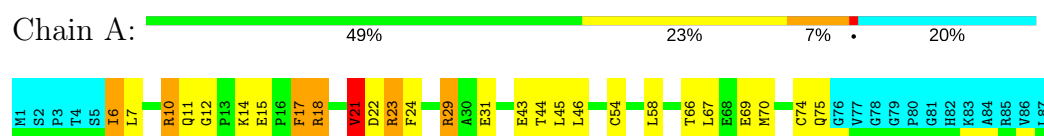
4.2.11 Score per residue for model 11

- Molecule 1: Gag-Pol polyprotein



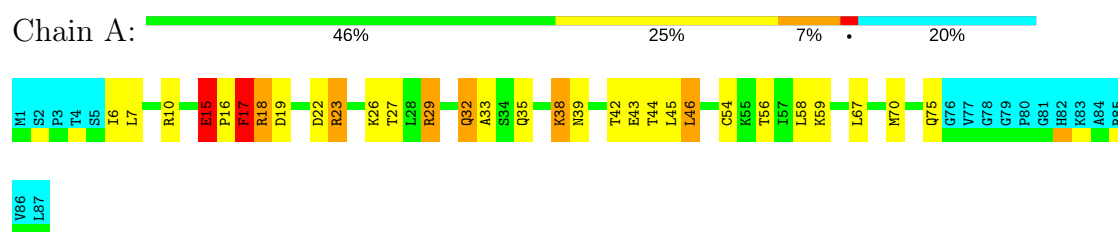
4.2.12 Score per residue for model 12

- Molecule 1: Gag-Pol polyprotein



4.2.13 Score per residue for model 13

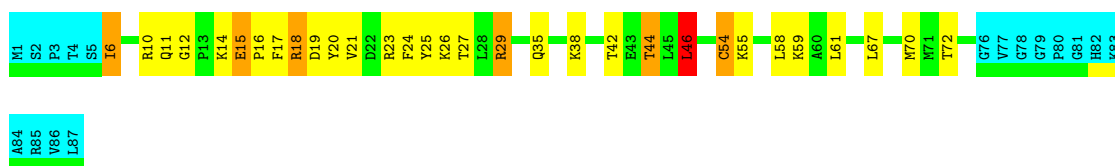
- Molecule 1: Gag-Pol polyprotein



4.2.14 Score per residue for model 14 (medoid)

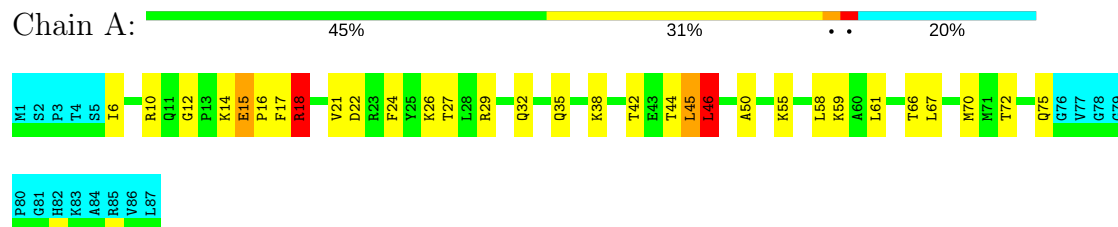
- Molecule 1: Gag-Pol polyprotein





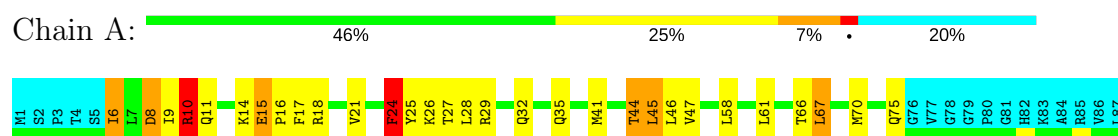
4.2.15 Score per residue for model 15

- Molecule 1: Gag-Pol polyprotein



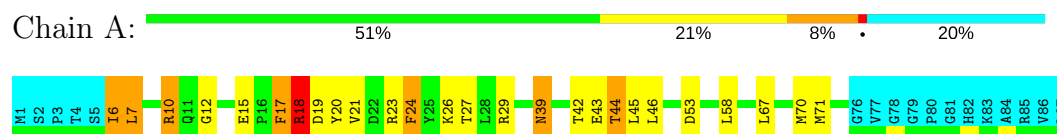
4.2.16 Score per residue for model 16

- Molecule 1: Gag-Pol polyprotein



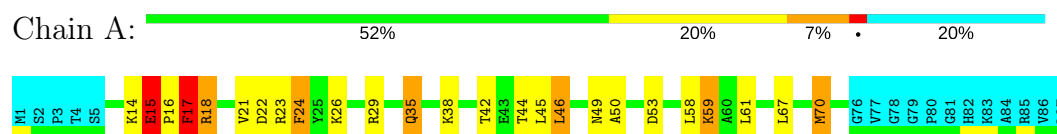
4.2.17 Score per residue for model 17

- Molecule 1: Gag-Pol polyprotein



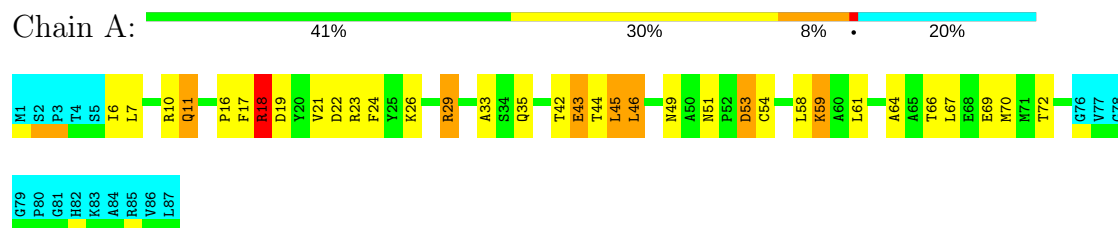
4.2.18 Score per residue for model 18

- Molecule 1: Gag-Pol polyprotein



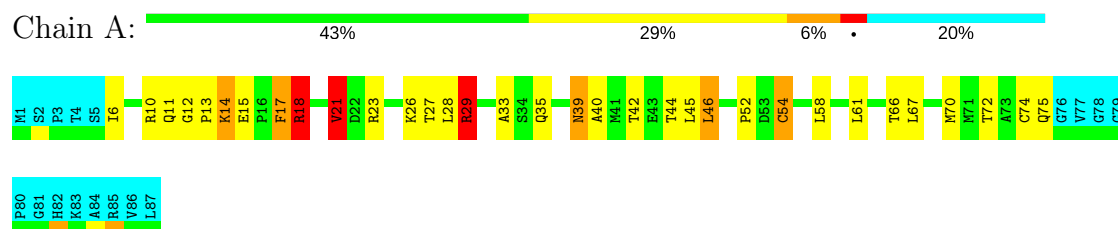
4.2.19 Score per residue for model 19

- Molecule 1: Gag-Pol polyprotein



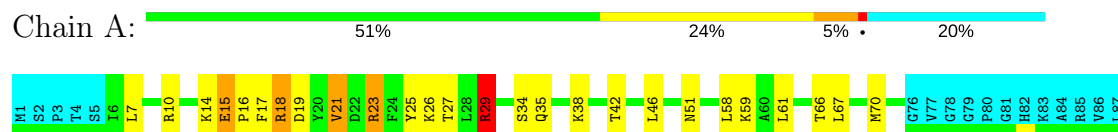
4.2.20 Score per residue for model 20

- Molecule 1: Gag-Pol polyprotein



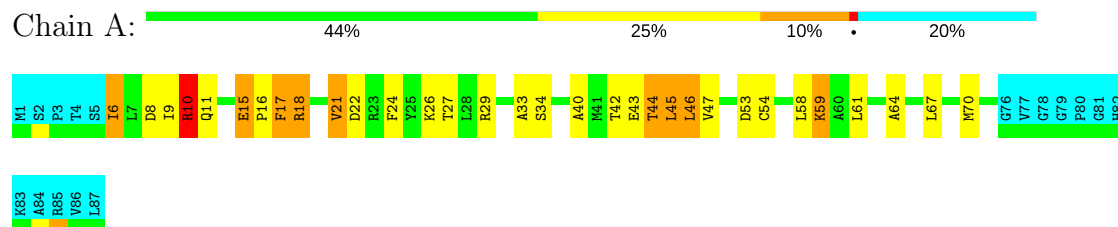
4.2.21 Score per residue for model 21

- Molecule 1: Gag-Pol polyprotein



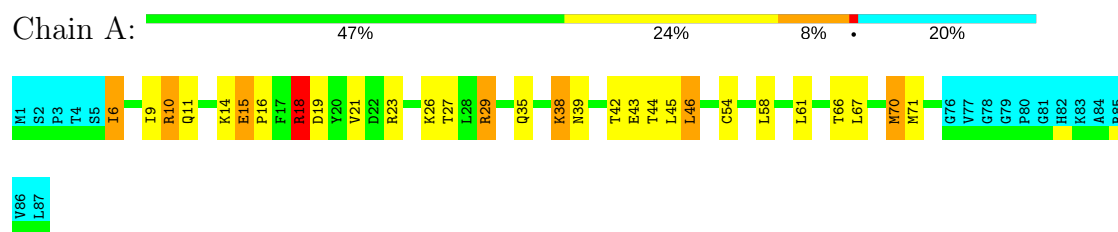
4.2.22 Score per residue for model 22

- Molecule 1: Gag-Pol polyprotein



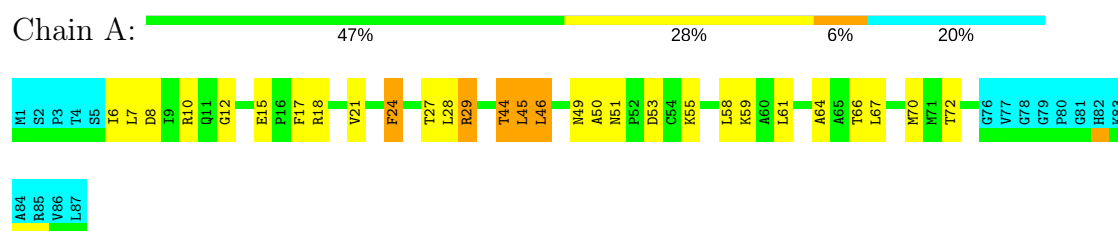
4.2.23 Score per residue for model 23

- Molecule 1: Gag-Pol polyprotein



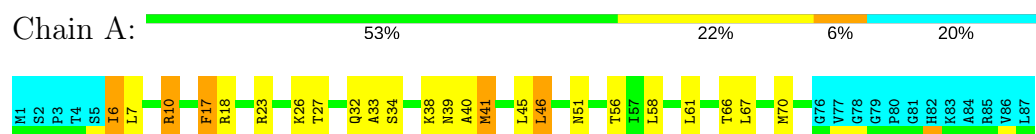
4.2.24 Score per residue for model 24

- Molecule 1: Gag-Pol polyprotein



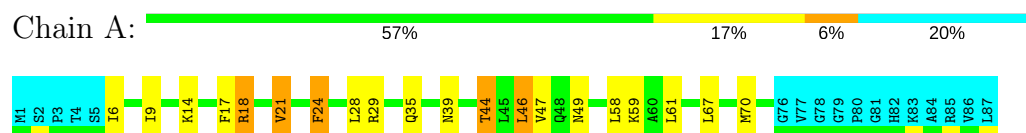
4.2.25 Score per residue for model 25

- Molecule 1: Gag-Pol polyprotein



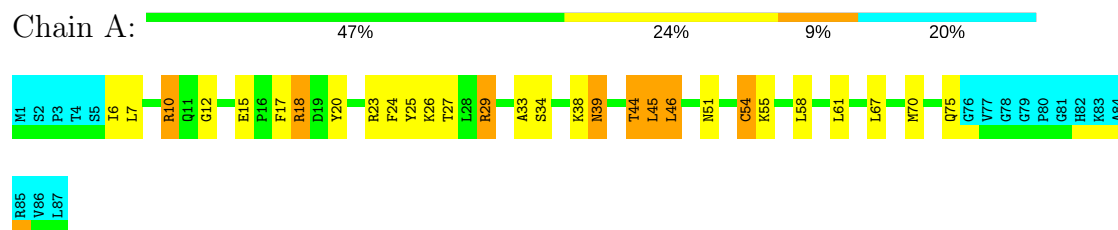
4.2.26 Score per residue for model 26

- Molecule 1: Gag-Pol polyprotein



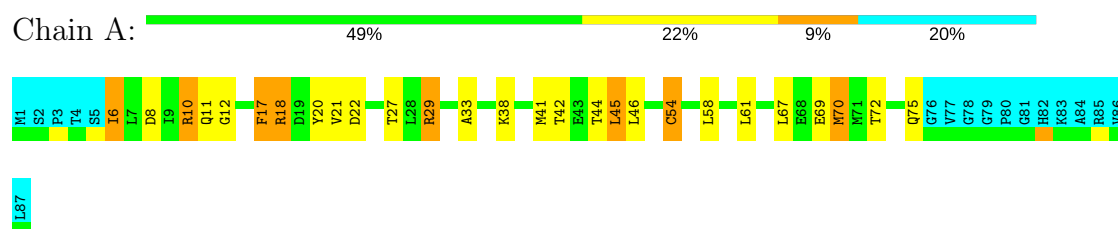
4.2.27 Score per residue for model 27

- Molecule 1: Gag-Pol polyprotein



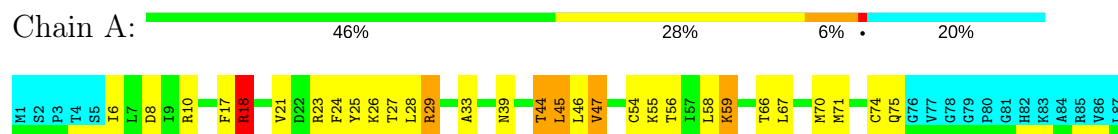
4.2.28 Score per residue for model 28

- Molecule 1: Gag-Pol polyprotein



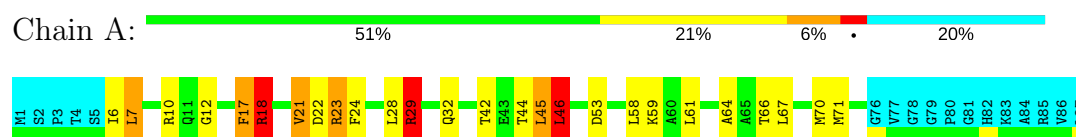
4.2.29 Score per residue for model 29

- Molecule 1: Gag-Pol polyprotein



4.2.30 Score per residue for model 30

- Molecule 1: Gag-Pol polyprotein



5 Refinement protocol and experimental data overview

The models were refined using the following method: *molecular dynamics*.

Of the 50 calculated structures, 30 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
CYANA	structure solution	2.1
AMBER	refinement	8

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 15137
Number of chemical shift lists	1
Total number of shifts	594
Number of shifts mapped to atoms	594
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	51%

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	0.70±0.01	0±0/559 (0.0±0.0%)	1.33±0.05	3±2/756 (0.4±0.2%)
All	All	0.70	0/16770 (0.0%)	1.33	89/22680 (0.4%)

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	0.9±0.7
All	All	0	27

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	23	ARG	NE-CZ-NH1	11.93	126.26	120.30	20	6
1	A	10	ARG	NE-CZ-NH1	10.28	125.44	120.30	4	10
1	A	18	ARG	NE-CZ-NH1	10.04	125.32	120.30	6	20
1	A	18	ARG	NE-CZ-NH2	-7.39	116.61	120.30	19	3
1	A	29	ARG	NE-CZ-NH1	7.13	123.87	120.30	21	8
1	A	21	VAL	CA-CB-CG1	6.84	121.15	110.90	20	4
1	A	18	ARG	CD-NE-CZ	6.49	132.68	123.60	19	2
1	A	10	ARG	NE-CZ-NH2	5.85	123.22	120.30	11	2
1	A	22	ASP	CB-CG-OD2	-5.83	113.05	118.30	12	2
1	A	46	LEU	CB-CG-CD2	5.82	120.89	111.00	3	10
1	A	24	PHE	CB-CG-CD2	-5.74	116.78	120.80	16	2
1	A	24	PHE	CB-CG-CD1	5.70	124.79	120.80	16	1
1	A	10	ARG	NH1-CZ-NH2	-5.56	113.29	119.40	16	2
1	A	7	LEU	C-N-CA	5.53	135.52	121.70	10	1
1	A	45	LEU	CB-CG-CD2	5.44	120.25	111.00	4	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	29	ARG	NE-CZ-NH2	-5.39	117.60	120.30	6	1
1	A	20	TYR	CB-CG-CD2	-5.38	117.77	121.00	6	2
1	A	9	ILE	C-N-CA	5.28	134.89	121.70	4	1
1	A	21	VAL	CA-CB-CG2	5.27	118.81	110.90	30	1
1	A	10	ARG	CD-NE-CZ	5.22	130.91	123.60	4	1
1	A	23	ARG	NE-CZ-NH2	-5.21	117.69	120.30	8	2
1	A	43	GLU	C-N-CA	5.21	134.72	121.70	22	3
1	A	50	ALA	CB-CA-C	5.14	117.81	110.10	24	1
1	A	46	LEU	CA-CB-CG	5.07	126.97	115.30	30	1
1	A	44	THR	CA-CB-CG2	5.02	119.43	112.40	26	1
1	A	10	ARG	N-CA-C	5.02	124.56	111.00	4	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	25	TYR	Sidechain	10
1	A	18	ARG	Sidechain	4
1	A	10	ARG	Sidechain	4
1	A	24	PHE	Sidechain	4
1	A	20	TYR	Sidechain	2
1	A	23	ARG	Sidechain	2
1	A	29	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	551	558	558	6±2
All	All	16530	16740	16740	189

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 6.

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:LEU:HD21	1:A:70:MET:HB2	0.78	1.54	16	27
1:A:58:LEU:HD21	1:A:70:MET:CB	0.70	2.15	10	27
1:A:21:VAL:HG23	1:A:46:LEU:HD23	0.64	1.68	8	3
1:A:21:VAL:HG12	1:A:46:LEU:HB2	0.63	1.71	24	1
1:A:24:PHE:CE1	1:A:44:THR:HG23	0.60	2.31	14	2
1:A:21:VAL:HG23	1:A:46:LEU:CD2	0.60	2.25	8	1
1:A:18:ARG:HA	1:A:21:VAL:HG12	0.60	1.74	9	11
1:A:21:VAL:HG13	1:A:46:LEU:HB2	0.59	1.74	30	1
1:A:47:VAL:HG22	1:A:71:MET:HA	0.58	1.74	29	1
1:A:17:PHE:CZ	1:A:58:LEU:HD22	0.57	2.33	12	11
1:A:58:LEU:HD21	1:A:70:MET:HB3	0.57	1.76	17	7
1:A:39:ASN:ND2	1:A:44:THR:HG22	0.57	2.15	2	2
1:A:39:ASN:HD21	1:A:44:THR:CG2	0.56	2.14	27	1
1:A:17:PHE:O	1:A:21:VAL:HG13	0.56	2.01	24	1
1:A:24:PHE:CZ	1:A:28:LEU:HD13	0.55	2.36	26	1
1:A:24:PHE:CZ	1:A:44:THR:HA	0.55	2.36	18	14
1:A:6:ILE:O	1:A:6:ILE:HD12	0.54	2.02	25	4
1:A:54:CYS:O	1:A:58:LEU:HD12	0.52	2.05	20	11
1:A:6:ILE:HD12	1:A:6:ILE:C	0.51	2.24	22	3
1:A:15:GLU:H	1:A:16:PRO:CD	0.51	2.18	13	2
1:A:39:ASN:HD22	1:A:44:THR:HG21	0.51	1.66	20	1
1:A:39:ASN:HD21	1:A:44:THR:HG22	0.51	1.66	27	1
1:A:54:CYS:SG	1:A:58:LEU:HD13	0.50	2.46	9	1
1:A:15:GLU:H	1:A:16:PRO:HD2	0.50	1.67	18	7
1:A:47:VAL:HG23	1:A:71:MET:HG2	0.50	1.82	5	1
1:A:36:GLU:CD	1:A:38:LYS:HE2	0.50	2.27	3	1
1:A:58:LEU:HD23	1:A:67:LEU:HG	0.49	1.84	16	1
1:A:16:PRO:HA	1:A:59:LYS:HE3	0.48	1.85	19	1
1:A:24:PHE:CZ	1:A:28:LEU:HD12	0.48	2.44	30	1
1:A:6:ILE:HD12	1:A:6:ILE:O	0.48	2.09	22	1
1:A:21:VAL:HG21	1:A:47:VAL:C	0.47	2.31	26	3
1:A:20:TYR:CD2	1:A:46:LEU:HD11	0.46	2.45	6	1
1:A:58:LEU:HD12	1:A:58:LEU:H	0.46	1.70	9	3
1:A:17:PHE:O	1:A:21:VAL:HG23	0.46	2.11	18	2
1:A:21:VAL:HG22	1:A:46:LEU:HG	0.46	1.88	18	1
1:A:21:VAL:HG23	1:A:46:LEU:HG	0.46	1.88	14	1
1:A:18:ARG:HA	1:A:21:VAL:HG22	0.45	1.87	2	1
1:A:28:LEU:HD13	1:A:28:LEU:C	0.45	2.31	11	1
1:A:39:ASN:HD22	1:A:44:THR:HG22	0.45	1.71	17	1
1:A:7:LEU:HD23	1:A:7:LEU:H	0.45	1.71	9	1
1:A:18:ARG:CD	1:A:18:ARG:C	0.45	2.85	20	1
1:A:6:ILE:HG21	1:A:24:PHE:CE1	0.45	2.47	12	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:24:PHE:CE1	1:A:44:THR:HA	0.45	2.47	5	1
1:A:21:VAL:HG23	1:A:46:LEU:HD13	0.44	1.88	23	1
1:A:58:LEU:HD11	1:A:70:MET:HB2	0.44	1.87	8	1
1:A:58:LEU:CD2	1:A:70:MET:HB2	0.44	2.42	30	1
1:A:55:LYS:HE3	1:A:55:LYS:HA	0.43	1.88	6	1
1:A:6:ILE:HG13	1:A:24:PHE:CD1	0.43	2.48	16	1
1:A:18:ARG:C	1:A:18:ARG:HD2	0.43	2.34	28	1
1:A:14:LYS:HE2	1:A:14:LYS:O	0.43	2.13	8	1
1:A:21:VAL:HG23	1:A:46:LEU:CG	0.43	2.44	8	1
1:A:39:ASN:HD21	1:A:47:VAL:HG22	0.43	1.74	9	1
1:A:15:GLU:N	1:A:16:PRO:CD	0.43	2.82	13	1
1:A:58:LEU:H	1:A:58:LEU:HD12	0.42	1.75	8	1
1:A:18:ARG:HD3	1:A:18:ARG:C	0.42	2.34	20	1
1:A:13:PRO:CD	1:A:45:LEU:HD13	0.42	2.45	20	1
1:A:17:PHE:C	1:A:21:VAL:HG23	0.42	2.35	18	1
1:A:18:ARG:HD3	1:A:19:ASP:N	0.42	2.30	17	1
1:A:39:ASN:ND2	1:A:47:VAL:HG22	0.41	2.30	9	1
1:A:16:PRO:HA	1:A:59:LYS:CE	0.41	2.45	19	1
1:A:28:LEU:HD11	1:A:32:GLN:HE21	0.41	1.75	30	1
1:A:21:VAL:HG22	1:A:46:LEU:CG	0.41	2.46	18	1
1:A:6:ILE:HG23	1:A:31:GLU:OE2	0.41	2.16	12	1
1:A:46:LEU:HD22	1:A:46:LEU:N	0.41	2.31	15	1
1:A:54:CYS:HB2	1:A:74:CYS:CB	0.41	2.45	29	1
1:A:18:ARG:HA	1:A:21:VAL:CG1	0.40	2.47	21	1
1:A:46:LEU:N	1:A:46:LEU:HD22	0.40	2.32	10	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	70/87 (80%)	43±3 (62±4%)	20±3 (28±5%)	7±2 (10±3%)	1	9
All	All	2100/2610 (80%)	1292 (62%)	588 (28%)	220 (10%)	1	9

All 35 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	15	GLU	20
1	A	44	THR	19
1	A	33	ALA	16
1	A	29	ARG	16
1	A	12	GLY	15
1	A	7	LEU	13
1	A	11	GLN	11
1	A	45	LEU	11
1	A	38	LYS	10
1	A	40	ALA	8
1	A	34	SER	7
1	A	49	ASN	7
1	A	64	ALA	6
1	A	39	ASN	6
1	A	14	LYS	6
1	A	10	ARG	6
1	A	59	LYS	5
1	A	17	PHE	4
1	A	6	ILE	4
1	A	35	GLN	3
1	A	50	ALA	3
1	A	32	GLN	3
1	A	43	GLU	3
1	A	51	ASN	3
1	A	75	GLN	2
1	A	18	ARG	2
1	A	41	MET	2
1	A	53	ASP	2
1	A	8	ASP	1
1	A	16	PRO	1
1	A	66	THR	1
1	A	52	PRO	1
1	A	13	PRO	1
1	A	9	ILE	1
1	A	48	GLN	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	60/72 (83%)	42±3 (70±6%)	18±3 (30±6%)	2	16
All	All	1800/2160 (83%)	1256 (70%)	544 (30%)	2	16

All 51 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	46	LEU	30
1	A	67	LEU	30
1	A	18	ARG	28
1	A	6	ILE	27
1	A	17	PHE	25
1	A	29	ARG	24
1	A	26	LYS	23
1	A	10	ARG	22
1	A	61	LEU	21
1	A	45	LEU	21
1	A	27	THR	19
1	A	42	THR	19
1	A	23	ARG	18
1	A	59	LYS	16
1	A	54	CYS	15
1	A	66	THR	14
1	A	35	GLN	13
1	A	75	GLN	13
1	A	72	THR	12
1	A	22	ASP	10
1	A	39	ASN	10
1	A	28	LEU	8
1	A	55	LYS	8
1	A	19	ASP	8
1	A	32	GLN	8
1	A	8	ASP	7
1	A	53	ASP	7
1	A	71	MET	7
1	A	38	LYS	7
1	A	44	THR	6
1	A	51	ASN	6
1	A	15	GLU	5
1	A	70	MET	5
1	A	43	GLU	5
1	A	14	LYS	5
1	A	11	GLN	4

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Mol	Chain	Res	Type	Models (Total)
1	A	7	LEU	4
1	A	69	GLU	4
1	A	21	VAL	4
1	A	56	THR	4
1	A	20	TYR	3
1	A	24	PHE	3
1	A	9	ILE	3
1	A	47	VAL	2
1	A	41	MET	2
1	A	74	CYS	2
1	A	37	VAL	2
1	A	49	ASN	2
1	A	34	SER	1
1	A	31	GLU	1
1	A	48	GLN	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.6 Ligand geometry ⓘ

There are no ligands in this entry.

6.7 Other polymers ⓘ

There are no such molecules in this entry.

6.8 Polymer linkage issues ⓘ

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 51% for the well-defined parts and 47% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 15137

Chemical shift list name: *assigned_chem_shift_list_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	594
Number of shifts mapped to atoms	594
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 [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	0	—	None (insufficient data)
$^{13}\text{C}_\beta$	0	—	None (insufficient data)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	68	-0.27 ± 0.41	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 51%, i.e. 442 atoms were assigned a chemical shift out of a possible 871. 0 out of 10 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	185/342 (54%)	125/136 (92%)	0/140 (0%)	60/66 (91%)
Sidechain	239/495 (48%)	239/290 (82%)	0/180 (0%)	0/25 (0%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	18/34 (53%)	18/18 (100%)	0/16 (0%)	0/0 (—%)
Overall	442/871 (51%)	382/444 (86%)	0/336 (0%)	60/91 (66%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 47%, i.e. 490 atoms were assigned a chemical shift out of a possible 1052. 0 out of 13 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	207/423 (49%)	139/168 (83%)	0/174 (0%)	68/81 (84%)
Sidechain	264/588 (45%)	264/346 (76%)	0/213 (0%)	0/29 (0%)
Aromatic	19/41 (46%)	19/22 (86%)	0/18 (0%)	0/1 (0%)
Overall	490/1052 (47%)	422/536 (79%)	0/405 (0%)	68/111 (61%)

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:

