



Full wwPDB NMR Structure Validation Report ⓘ

May 28, 2020 – 11:07 pm BST

PDB ID : 2LJZ
Title : Structure of the C-terminal domain of HPV16 E6 oncoprotein
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Deposited on : 2011-09-30

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 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 : 20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : 2.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

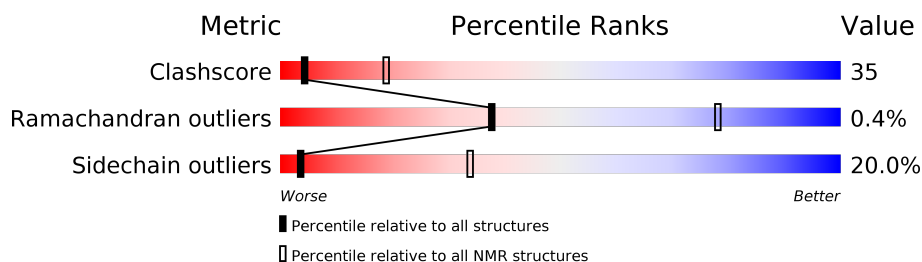
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 74%.

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	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

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	75	

2 Ensemble composition and analysis

This entry contains 20 models. Model 19 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:81-A:139 (59)	0.18	19

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1244 atoms, of which 623 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Protein E6.

Mol	Chain	Residues	Atoms						Trace
1	A	75	Total	C	H	N	O	S	0
			1243	377	623	124	113	6	

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	77	GLY	-	EXPRESSION TAG	UNP P03126
A	78	ALA	-	EXPRESSION TAG	UNP P03126
A	79	MET	-	EXPRESSION TAG	UNP P03126
A	80	SER	CYS	ENGINEERED MUTATION	UNP P03126
A	97	SER	CYS	ENGINEERED MUTATION	UNP P03126
A	111	SER	CYS	ENGINEERED MUTATION	UNP P03126
A	140	SER	CYS	ENGINEERED MUTATION	UNP P03126

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

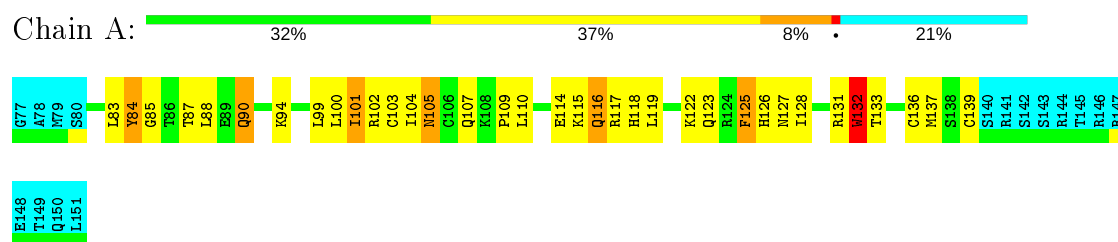
Mol	Chain	Residues	Atoms	
2	A	1	Total	Zn
			1	1

4 Residue-property plots [i](#)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA 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: Protein E6

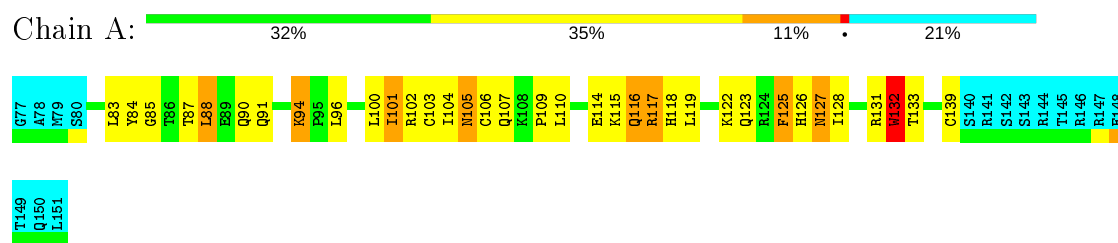


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

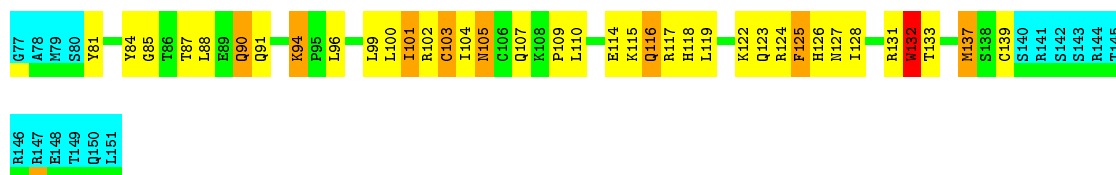
- Molecule 1: Protein E6



4.2.2 Score per residue for model 2

- Molecule 1: Protein E6

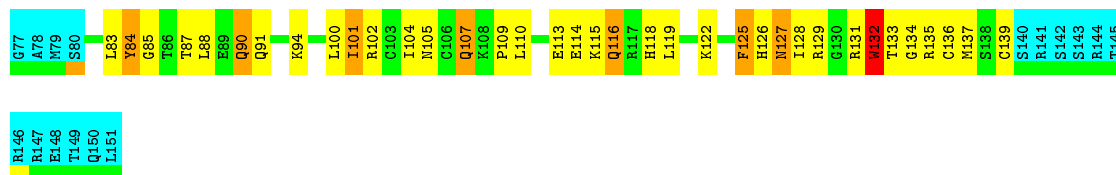




4.2.3 Score per residue for model 3

- Molecule 1: Protein E6

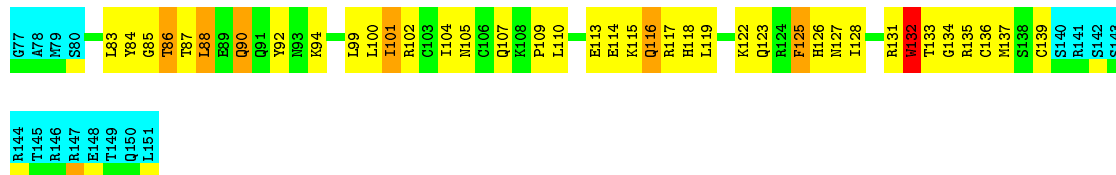
Chain A: 31% 37% 9% 21%



4.2.4 Score per residue for model 4

- Molecule 1: Protein E6

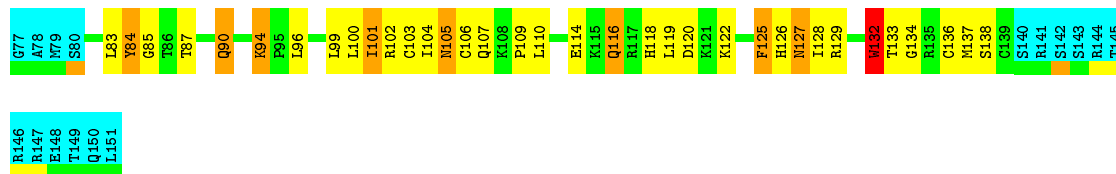
Chain A: 27% 43% 8% 21%



4.2.5 Score per residue for model 5

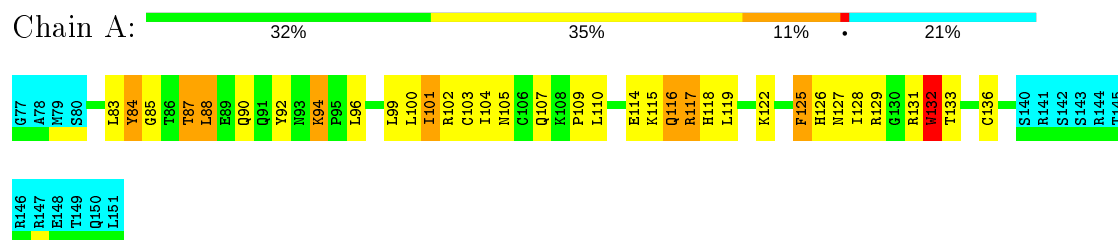
- Molecule 1: Protein E6

Chain A: 32% 35% 11% 21%



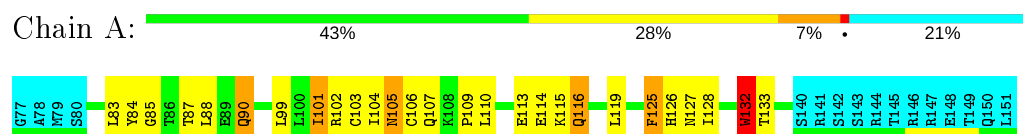
4.2.6 Score per residue for model 6

- Molecule 1: Protein E6



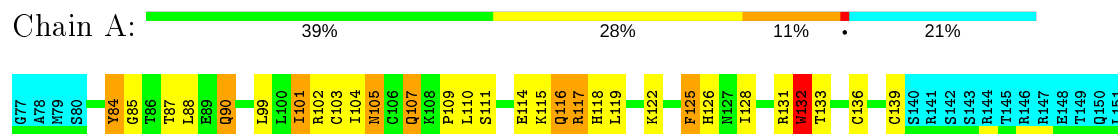
4.2.7 Score per residue for model 7

- Molecule 1: Protein E6



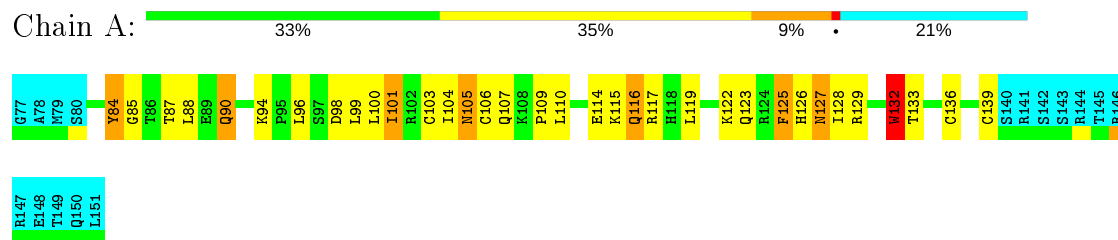
4.2.8 Score per residue for model 8

- Molecule 1: Protein E6



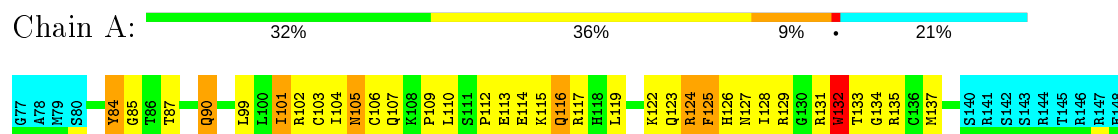
4.2.9 Score per residue for model 9

- Molecule 1: Protein E6



4.2.10 Score per residue for model 10

- Molecule 1: Protein E6



T149
Q150
L151

4.2.11 Score per residue for model 11

- Molecule 1: Protein E6

Chain A: 36% 28% 13% • 21%

G77 A78 M79 S80 L83 Y84 G85 T86 T87 L88 L89 Q90 R94 L99 L100 I101 R102 C103 I104 M105 M106 Q107 K108 P109 L110 E114 K115 Q116 L117 L119 K122 F125 H126 M127 I128 R131 R132 T133 G134 R135 C136 C139 S140 R141 S142 S143 S144 T145 T146 R147 R148 R149 Q150 L151

Q150
L151

4.2.12 Score per residue for model 12

- Molecule 1: Protein E6

Chain A: 36% 33% 8% • 21%

G77 A78 M79 S80 L83 Y84 G85 T86 T87 L88 L89 Q90 R94 L99 L100 I101 R102 C103 I104 M105 M106 Q107 K108 P109 L110 E114 K115 Q116 L117 L119 K122 F125 H126 M127 I128 R131 R132 T133 M137 S140 R141 S142 S143 T144 T145 T146 R147 R148 R149 Q150 L151

4.2.13 Score per residue for model 13

- Molecule 1: Protein E6

Chain A: 31% 36% 11% • 21%

G77 A78 M79 S80 L83 Y84 G85 T86 T87 L88 L89 Q90 R94 R95 L99 L100 S97 D98 L99 L100 I101 M105 M106 Q107 M108 P109 L110 E114 K115 Q116 L117 L119 K122 Q123 R124 F125 H126 M127 I128 R131 R132 T133 C136 C139 S140 R141 S142 S143 R144 T145 T146 R147 R148 R149 Q150 L151

R147
E148
T149
Q150
L151

4.2.14 Score per residue for model 14

- Molecule 1: Protein E6

Chain A: 31% 37% 9% • 21%

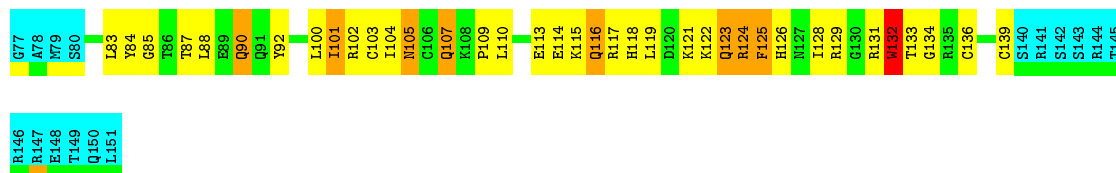
G77 A78 M79 S80 L83 Y84 G85 T86 T87 L88 L89 Q90 R94 L99 L100 I101 R102 C103 I104 M105 M106 Q107 K108 P109 L110 S111 P112 E113 E114 K115 Q116 L117 L119 K122 R123 F125 H126 M127 I128 R131 R132 T133 G134 R135 S140 R141 S142 S143 R144 T145 T146 R147 R148 R149 Q150 L151

Q150
L151

4.2.15 Score per residue for model 15

- Molecule 1: Protein E6

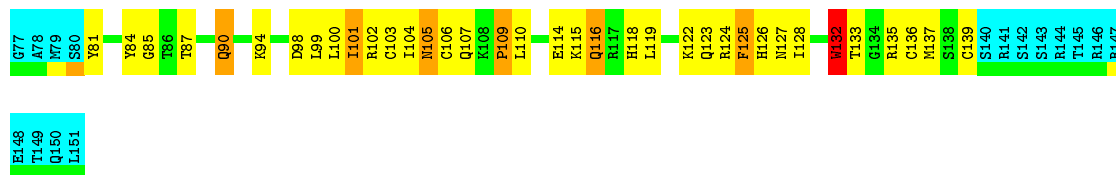
Chain A: 29% 37% 11% 21%



4.2.16 Score per residue for model 16

- Molecule 1: Protein E6

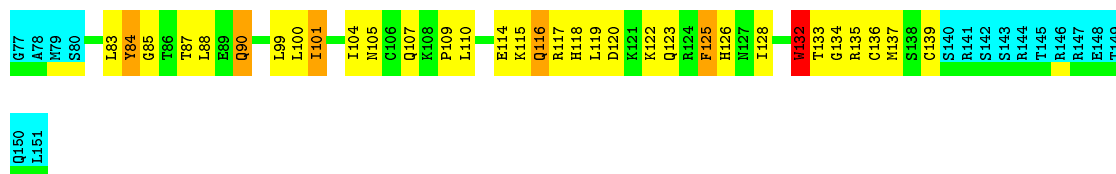
Chain A: 31% 39% 8% 21%



4.2.17 Score per residue for model 17

- Molecule 1: Protein E6

Chain A: 35% 36% 7% 21%



4.2.18 Score per residue for model 18

- Molecule 1: Protein E6

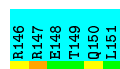
Chain A: 40% 28% 9% 21%



4.2.19 Score per residue for model 19 (medoid)

- Molecule 1: Protein E6

Chain A: 32% 36% 9% 21%



4.2.20 Score per residue for model 20

- Molecule 1: Protein E6

Chain A: 32% 35% 11% 21%



5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 20 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
X-PLOR NIH	structure solution	
X-PLOR NIH	refinement	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 6 of this report.

Chemical shift file(s)	input_cs.cif
Number of chemical shift lists	1
Total number of shifts	788
Number of shifts mapped to atoms	788
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	74%

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

COVALENT-GEOMETRY INFOmissingINFO

5.1 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	494	493	493	35±3
All	All	9900	9860	9860	697

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 35.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:85:GLY:HA2	1:A:125:PHE:CZ	0.86	2.05	3	20
1:A:104:ILE:HA	1:A:133:THR:OG1	0.76	1.80	13	20
1:A:110:LEU:HB3	1:A:114:GLU:HB2	0.75	1.59	17	20
1:A:87:THR:O	1:A:90:GLN:HG3	0.67	1.89	6	3
1:A:101:ILE:HG12	1:A:132:TRP:CD2	0.67	2.24	6	20
1:A:128:ILE:HD13	1:A:132:TRP:HA	0.65	1.69	17	20
1:A:87:THR:O	1:A:90:GLN:HB3	0.65	1.92	17	16
1:A:134:GLY:O	1:A:135:ARG:HG3	0.64	1.92	3	5
1:A:117:ARG:NH2	1:A:137:MET:HG3	0.64	2.08	2	1
1:A:101:ILE:HG12	1:A:132:TRP:CE2	0.63	2.28	20	20
1:A:114:GLU:OE1	1:A:137:MET:HG3	0.63	1.94	17	1
1:A:104:ILE:HB	1:A:134:GLY:O	0.61	1.95	3	7
1:A:121:LYS:HB2	1:A:123:GLN:OE1	0.60	1.95	15	1
1:A:118:HIS:O	1:A:122:LYS:HA	0.60	1.95	12	15
1:A:128:ILE:HD11	1:A:133:THR:HG22	0.60	1.74	17	20
1:A:101:ILE:HG23	1:A:132:TRP:CH2	0.59	2.32	11	20
1:A:105:ASN:ND2	1:A:139:CYS:SG	0.59	2.75	4	4
1:A:126:HIS:O	1:A:132:TRP:HB2	0.58	1.99	13	18
1:A:115:LYS:O	1:A:119:LEU:HG	0.56	2.00	14	19
1:A:103:CYS:HB3	1:A:106:CYS:HB2	0.56	1.77	19	10
1:A:102:ARG:O	1:A:133:THR:HA	0.56	2.00	6	17
1:A:101:ILE:O	1:A:109:PRO:HA	0.55	2.02	1	20
1:A:84:TYR:CD2	1:A:122:LYS:HB3	0.54	2.37	8	13
1:A:116:GLN:CA	1:A:116:GLN:HE21	0.54	2.16	20	9
1:A:114:GLU:OE2	1:A:136:CYS:HB2	0.54	2.02	18	10
1:A:116:GLN:HE21	1:A:116:GLN:CA	0.54	2.16	2	9
1:A:132:TRP:HE3	1:A:133:THR:N	0.53	2.02	5	20
1:A:105:ASN:H	1:A:105:ASN:ND2	0.53	2.00	8	6
1:A:129:ARG:O	1:A:129:ARG:HD2	0.53	2.03	15	1
1:A:105:ASN:ND2	1:A:105:ASN:H	0.52	2.00	13	5
1:A:103:CYS:CB	1:A:106:CYS:HB2	0.52	2.34	19	7
1:A:101:ILE:HG23	1:A:132:TRP:CZ3	0.52	2.40	2	20
1:A:105:ASN:ND2	1:A:139:CYS:HB3	0.52	2.19	8	9
1:A:107:GLN:HE21	1:A:107:GLN:N	0.51	2.03	8	1
1:A:118:HIS:HD1	1:A:123:GLN:HB2	0.51	1.65	20	4
1:A:85:GLY:CA	1:A:125:PHE:CZ	0.51	2.94	19	13
1:A:99:LEU:HD23	1:A:101:ILE:HD11	0.50	1.84	12	17
1:A:105:ASN:HD21	1:A:139:CYS:CB	0.50	2.20	4	3
1:A:128:ILE:HB	1:A:131:ARG:O	0.50	2.07	3	8
1:A:102:ARG:HD3	1:A:107:GLN:O	0.50	2.07	3	4
1:A:94:LYS:HG3	1:A:98:ASP:HB3	0.50	1.83	9	3
1:A:103:CYS:SG	1:A:105:ASN:ND2	0.49	2.86	9	14

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:113:GLU:HG3	1:A:114:GLU:OE1	0.49	2.06	3	6
1:A:114:GLU:HG2	1:A:137:MET:HG3	0.49	1.85	16	3
1:A:131:ARG:NH1	1:A:133:THR:HB	0.48	2.23	2	5
1:A:125:PHE:CE2	1:A:132:TRP:CZ2	0.48	3.01	3	17
1:A:114:GLU:CD	1:A:136:CYS:HB2	0.48	2.29	11	2
1:A:116:GLN:HE21	1:A:116:GLN:N	0.48	2.07	7	11
1:A:125:PHE:CD2	1:A:132:TRP:CZ2	0.48	3.02	1	16
1:A:100:LEU:O	1:A:131:ARG:HA	0.48	2.08	2	4
1:A:117:ARG:O	1:A:123:GLN:HG2	0.48	2.09	2	5
1:A:84:TYR:HB3	1:A:86:THR:HG22	0.47	1.85	4	1
1:A:114:GLU:O	1:A:117:ARG:HB2	0.47	2.09	6	1
1:A:117:ARG:HG2	1:A:123:GLN:CG	0.47	2.39	4	1
1:A:110:LEU:HD22	1:A:136:CYS:HB3	0.47	1.86	15	3
1:A:116:GLN:N	1:A:116:GLN:HE21	0.47	2.08	20	7
1:A:87:THR:HB	1:A:90:GLN:NE2	0.47	2.25	20	4
1:A:84:TYR:CE2	1:A:122:LYS:HB3	0.47	2.44	12	5
1:A:112:PRO:O	1:A:116:GLN:HB2	0.46	2.09	19	3
1:A:123:GLN:O	1:A:124:ARG:HG2	0.46	2.11	2	1
1:A:83:LEU:HD23	1:A:88:LEU:HD22	0.46	1.86	3	11
1:A:116:GLN:NE2	1:A:119:LEU:HD12	0.46	2.25	11	10
1:A:86:THR:HG21	1:A:122:LYS:HE3	0.45	1.88	4	1
1:A:116:GLN:O	1:A:120:ASP:HB3	0.45	2.12	5	1
1:A:132:TRP:C	1:A:132:TRP:HE3	0.45	2.15	2	8
1:A:94:LYS:HD2	1:A:99:LEU:CD1	0.45	2.42	20	3
1:A:94:LYS:HG3	1:A:94:LYS:O	0.45	2.12	5	1
1:A:119:LEU:O	1:A:122:LYS:HG2	0.44	2.13	2	2
1:A:132:TRP:HE3	1:A:132:TRP:C	0.44	2.16	3	4
1:A:128:ILE:HD13	1:A:132:TRP:CA	0.44	2.42	17	5
1:A:127:ASN:C	1:A:128:ILE:HD12	0.44	2.33	5	3
1:A:128:ILE:HD12	1:A:128:ILE:N	0.44	2.28	9	4
1:A:117:ARG:HG2	1:A:123:GLN:HG3	0.44	1.90	1	2
1:A:117:ARG:HB2	1:A:117:ARG:CZ	0.44	2.43	8	2
1:A:88:LEU:HB2	1:A:96:LEU:HD11	0.44	1.88	9	2
1:A:102:ARG:HG2	1:A:109:PRO:CA	0.44	2.43	15	6
1:A:128:ILE:CD1	1:A:133:THR:HG22	0.44	2.43	9	5
1:A:85:GLY:HA2	1:A:125:PHE:HZ	0.43	1.72	14	1
1:A:83:LEU:HD23	1:A:88:LEU:CD2	0.43	2.43	13	4
1:A:83:LEU:HG	1:A:84:TYR:N	0.43	2.29	5	1
1:A:125:PHE:CE2	1:A:132:TRP:CE2	0.43	3.07	16	3
1:A:118:HIS:ND1	1:A:123:GLN:HB2	0.43	2.28	20	3
1:A:132:TRP:CE3	1:A:133:THR:N	0.43	2.85	1	12

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:99:LEU:HD23	1:A:101:ILE:CD1	0.43	2.44	17	4
1:A:110:LEU:CD2	1:A:136:CYS:HB3	0.43	2.43	13	3
1:A:96:LEU:CD1	1:A:119:LEU:HD21	0.43	2.43	9	1
1:A:117:ARG:O	1:A:123:GLN:HG3	0.42	2.13	20	1
1:A:116:GLN:O	1:A:120:ASP:HB2	0.42	2.15	17	1
1:A:105:ASN:HD21	1:A:139:CYS:HB3	0.42	1.75	18	2
1:A:118:HIS:ND1	1:A:123:GLN:HB3	0.42	2.30	16	1
1:A:105:ASN:ND2	1:A:139:CYS:CB	0.41	2.83	18	1
1:A:104:ILE:HG22	1:A:135:ARG:HG3	0.41	1.91	17	1
1:A:104:ILE:HG13	1:A:133:THR:OG1	0.41	2.15	3	2
1:A:96:LEU:HD13	1:A:119:LEU:HD21	0.41	1.93	5	1
1:A:92:TYR:C	1:A:94:LYS:H	0.41	2.18	6	2
1:A:96:LEU:HB3	1:A:115:LYS:HD3	0.41	1.91	6	3
1:A:104:ILE:HA	1:A:133:THR:HG1	0.41	1.75	6	1
1:A:110:LEU:HB3	1:A:114:GLU:CB	0.41	2.39	17	1
1:A:128:ILE:CD1	1:A:132:TRP:HA	0.41	2.45	12	1
1:A:105:ASN:N	1:A:105:ASN:ND2	0.41	2.67	8	1
1:A:111:SER:HB2	1:A:113:GLU:HG2	0.40	1.93	14	1
1:A:124:ARG:HD2	1:A:124:ARG:H	0.40	1.76	10	1
1:A:94:LYS:O	1:A:94:LYS:HG3	0.40	2.15	1	2
1:A:102:ARG:HA	1:A:109:PRO:HA	0.40	1.94	1	2
1:A:110:LEU:O	1:A:115:LYS:HE3	0.40	2.17	6	1
1:A:94:LYS:HE3	1:A:99:LEU:HA	0.40	1.93	12	1

5.2 Torsion angles ⓘ

5.2.1 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	59/75 (79%)	54±2 (91±3%)	5±1 (9±2%)	0±1 (0±1%)	38	78
All	All	1180/1500 (79%)	1074 (91%)	101 (9%)	5 (0%)	38	78

All 4 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	129	ARG	2
1	A	86	THR	1
1	A	123	GLN	1
1	A	124	ARG	1

5.2.2 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	56/70 (80%)	45±2 (80±3%)	11±2 (20±3%)	4	34
All	All	1120/1400 (80%)	896 (80%)	224 (20%)	4	34

All 23 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	132	TRP	20
1	A	125	PHE	20
1	A	116	GLN	20
1	A	101	ILE	20
1	A	107	GLN	20
1	A	105	ASN	17
1	A	90	GLN	17
1	A	84	TYR	16
1	A	127	ASN	15
1	A	100	LEU	13
1	A	94	LYS	9
1	A	88	LEU	7
1	A	124	ARG	4
1	A	135	ARG	4
1	A	91	GLN	4
1	A	137	MET	4
1	A	103	CYS	4
1	A	129	ARG	3
1	A	117	ARG	2
1	A	87	THR	2
1	A	123	GLN	1
1	A	111	SER	1

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Mol	Chain	Res	Type	Models (Total)
1	A	109	PRO	1

5.2.3 RNA [i](#)

There are no RNA molecules in this entry.

5.3 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.4 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.5 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

5.6 Other polymers [i](#)

There are no such molecules in this entry.

5.7 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Chemical shift validation [i](#)

The completeness of assignment taking into account all chemical shift lists is 74% for the well-defined parts and 71% for the entire structure.

6.1 Chemical shift list 1

File name: input_cs.cif

Chemical shift list name: *E6C*

6.1.1 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	788
Number of shifts mapped to atoms	788
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	2

6.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	73	-0.75 ± 0.17	Should be applied
$^{13}\text{C}_\beta$	69	0.41 ± 0.15	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	69	0.49 ± 0.50	None needed (< 0.5 ppm)

6.1.3 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 74%, i.e. 597 atoms were assigned a chemical shift out of a possible 803. 7 out of 7 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	229/289 (79%)	115/115 (100%)	58/118 (49%)	56/56 (100%)
Sidechain	348/455 (76%)	217/273 (79%)	123/151 (81%)	8/31 (26%)

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	Total	¹H	¹³C	¹⁵N
Aromatic	20/59 (34%)	19/31 (61%)	0/25 (0%)	1/3 (33%)
Overall	597/803 (74%)	351/419 (84%)	181/294 (62%)	65/90 (72%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 71%, i.e. 715 atoms were assigned a chemical shift out of a possible 1011. 8 out of 8 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	283/369 (77%)	141/147 (96%)	73/150 (49%)	69/72 (96%)
Sidechain	412/583 (71%)	254/350 (73%)	149/189 (79%)	9/44 (20%)
Aromatic	20/59 (34%)	19/31 (61%)	0/25 (0%)	1/3 (33%)
Overall	715/1011 (71%)	414/528 (78%)	222/364 (61%)	79/119 (66%)

6.1.4 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	127	ASN	HB3	-0.30	4.41 – 1.11	-9.3
1	A	127	ASN	HD22	3.86	9.59 – 4.69	-6.7

6.1.5 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:

