



Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 07:06 PM BST

PDB ID : 2CQK
Title : Solution structure of the La domain of c-Mpl binding protein
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Deposited on : 2005-05-20

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

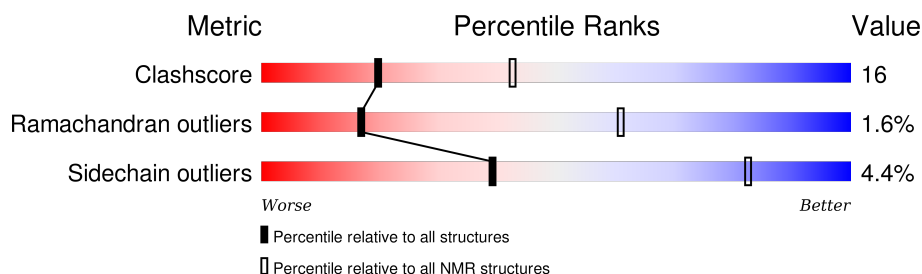
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

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	101	<div> <div></div> <div>55%</div> <div>23%</div> <div>22%</div> </div>

2 Ensemble composition and analysis

This entry contains 20 models. Model 11 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:47-A:125 (79)	0.22	11

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1568 atoms, of which 786 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called C-Mpl binding protein.

Mol	Chain	Residues	Atoms						Trace
1	A	101	Total	C	H	N	O	S	0
			1568	486	786	131	159	6	

There are 13 discrepancies between the modelled and reference sequences:

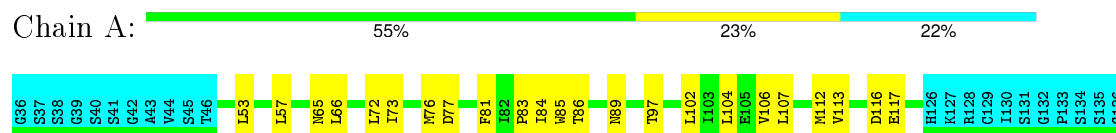
Chain	Residue	Modelled	Actual	Comment	Reference
A	36	GLY	-	CLONING ARTIFACT	UNP Q71RC2
A	37	SER	-	CLONING ARTIFACT	UNP Q71RC2
A	38	SER	-	CLONING ARTIFACT	UNP Q71RC2
A	39	GLY	-	CLONING ARTIFACT	UNP Q71RC2
A	40	SER	-	CLONING ARTIFACT	UNP Q71RC2
A	41	SER	-	CLONING ARTIFACT	UNP Q71RC2
A	42	GLY	-	CLONING ARTIFACT	UNP Q71RC2
A	131	SER	-	CLONING ARTIFACT	UNP Q71RC2
A	132	GLY	-	CLONING ARTIFACT	UNP Q71RC2
A	133	PRO	-	CLONING ARTIFACT	UNP Q71RC2
A	134	SER	-	CLONING ARTIFACT	UNP Q71RC2
A	135	SER	-	CLONING ARTIFACT	UNP Q71RC2
A	136	GLY	-	CLONING ARTIFACT	UNP Q71RC2

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: C-Mpl binding protein

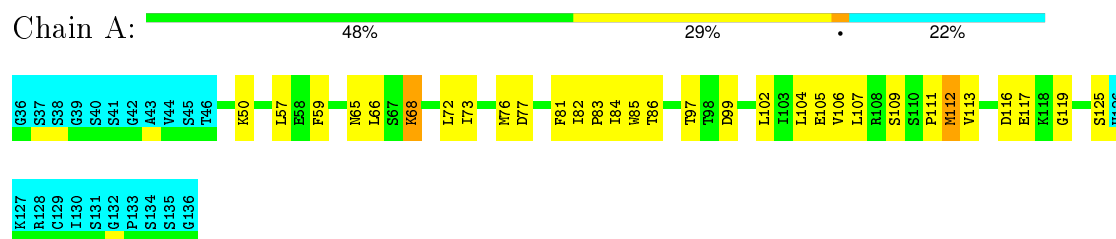


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

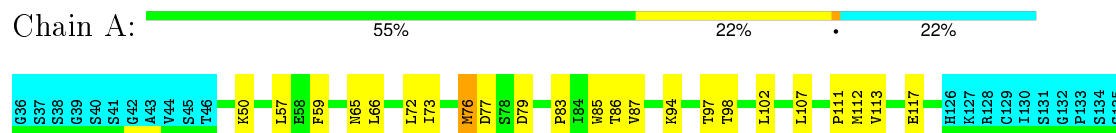
4.2.1 Score per residue for model 1

- Molecule 1: C-Mpl binding protein



4.2.2 Score per residue for model 2

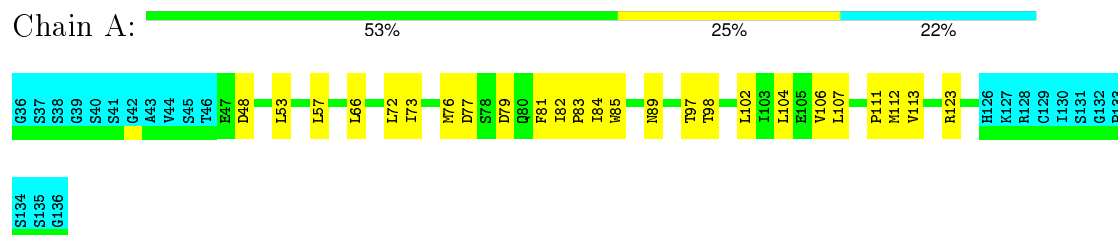
- Molecule 1: C-Mpl binding protein



G136

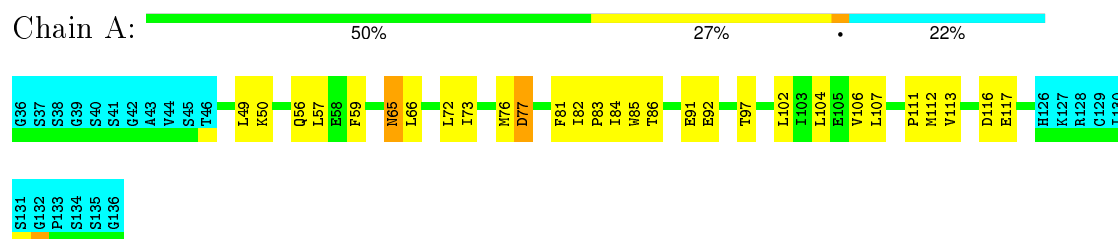
4.2.3 Score per residue for model 3

- Molecule 1: C-Mpl binding protein

S134
S135
G136

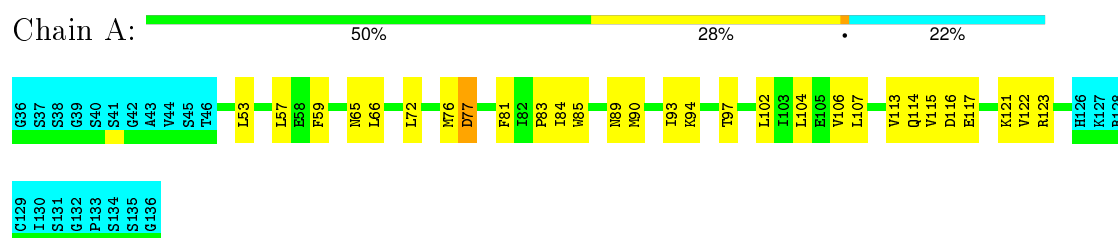
4.2.4 Score per residue for model 4

- Molecule 1: C-Mpl binding protein

S131
G132
P133
S134
S135
G136

4.2.5 Score per residue for model 5

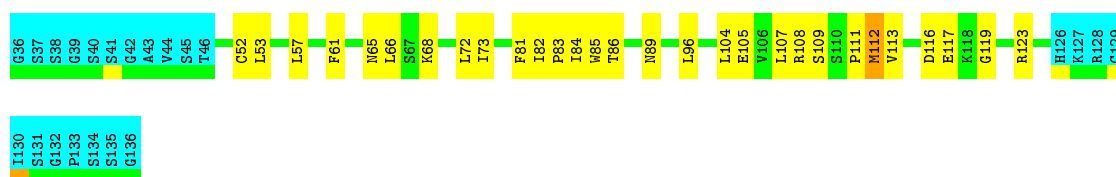
- Molecule 1: C-Mpl binding protein

G129
I130
S131
G132
P133
S134
S135
G136

4.2.6 Score per residue for model 6

- Molecule 1: C-Mpl binding protein

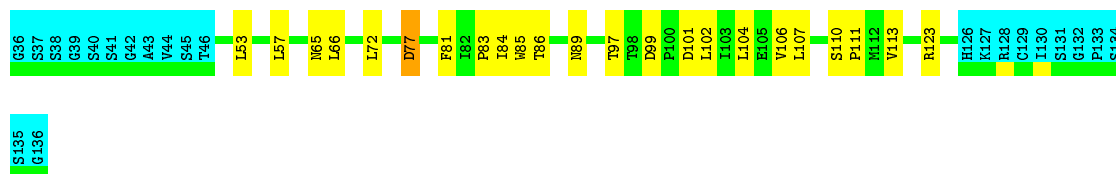




4.2.7 Score per residue for model 7

- Molecule 1: C-Mpl binding protein

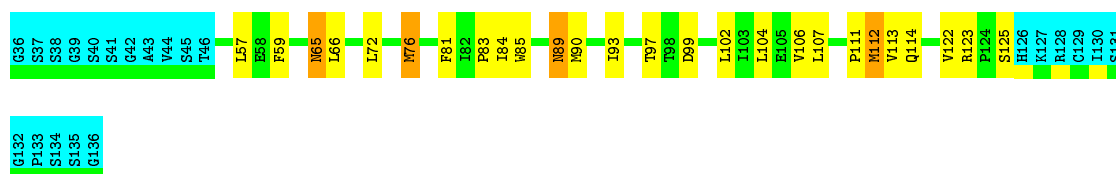
Chain A: 55% 22% 22%



4.2.8 Score per residue for model 8

- Molecule 1: C-Mpl binding protein

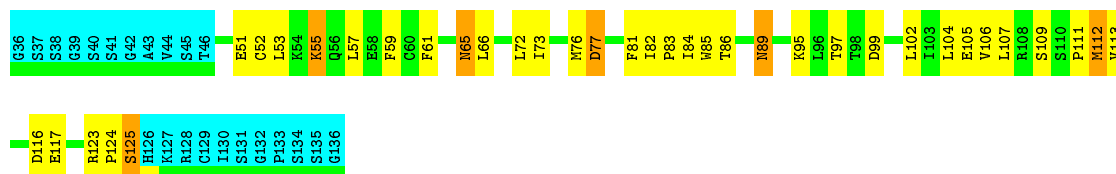
Chain A: 52% 22% 22%



4.2.9 Score per residue for model 9

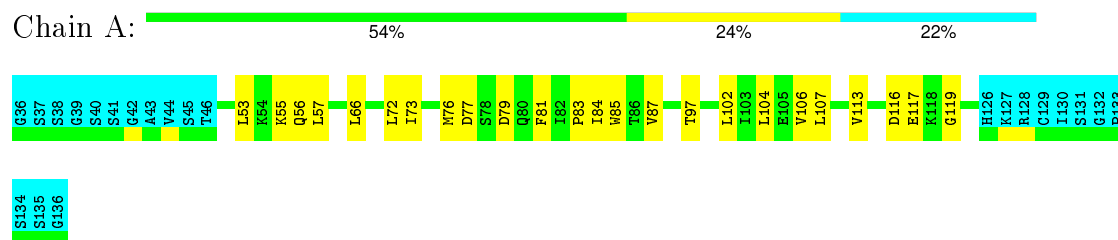
- Molecule 1: C-Mpl binding protein

Chain A: 42% 31% 6% 22%



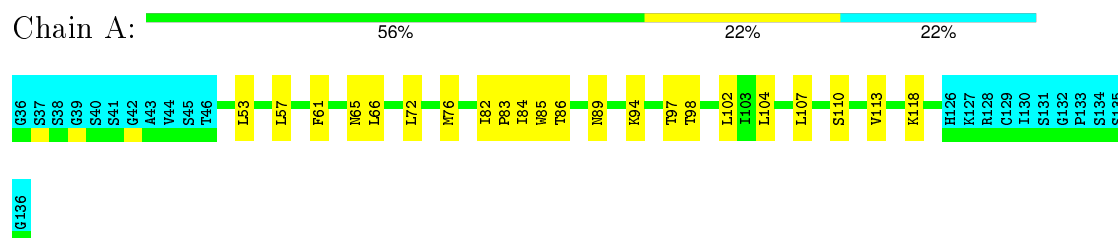
4.2.10 Score per residue for model 10

- Molecule 1: C-Mpl binding protein



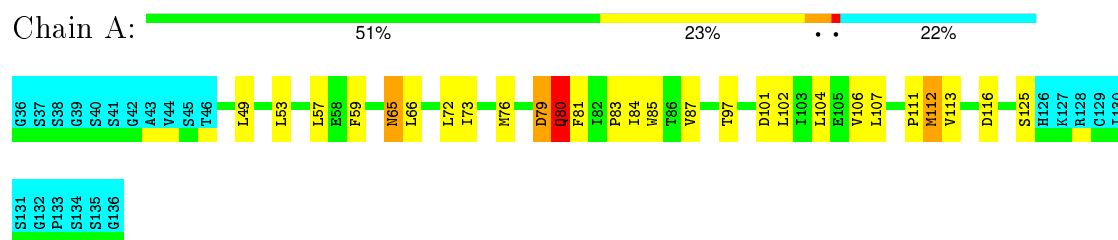
4.2.11 Score per residue for model 11 (medoid)

- Molecule 1: C-Mpl binding protein



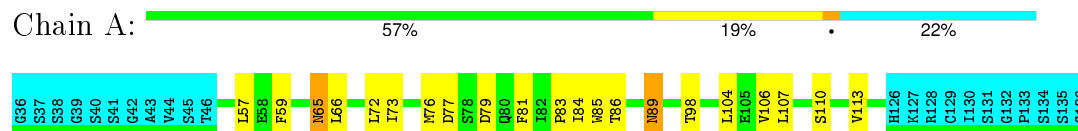
4.2.12 Score per residue for model 12

- Molecule 1: C-Mpl binding protein



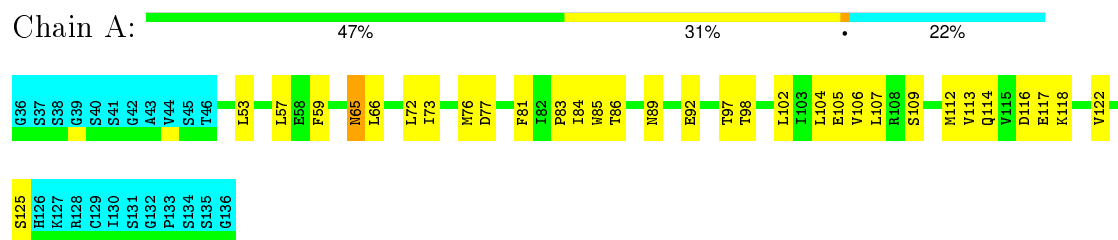
4.2.13 Score per residue for model 13

- Molecule 1: C-Mpl binding protein



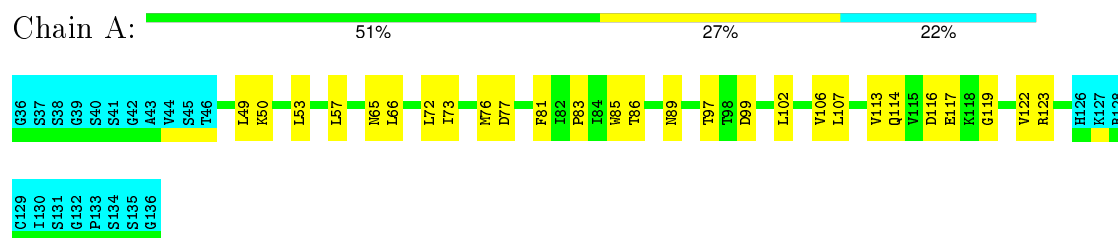
4.2.14 Score per residue for model 14

- Molecule 1: C-Mpl binding protein



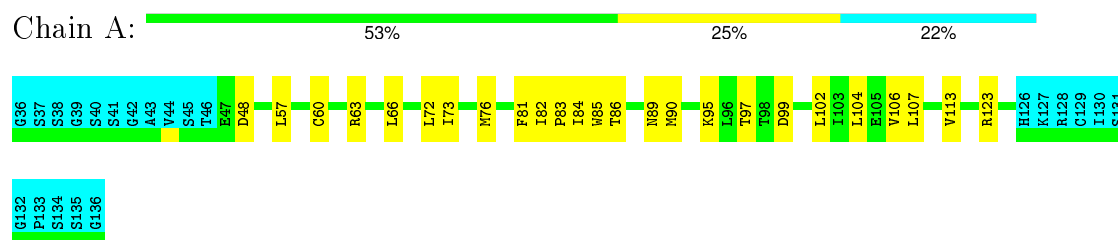
4.2.15 Score per residue for model 15

- Molecule 1: C-Mpl binding protein



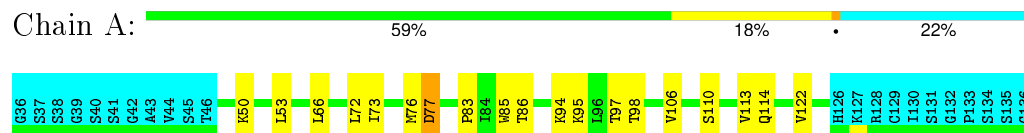
4.2.16 Score per residue for model 16

- Molecule 1: C-Mpl binding protein



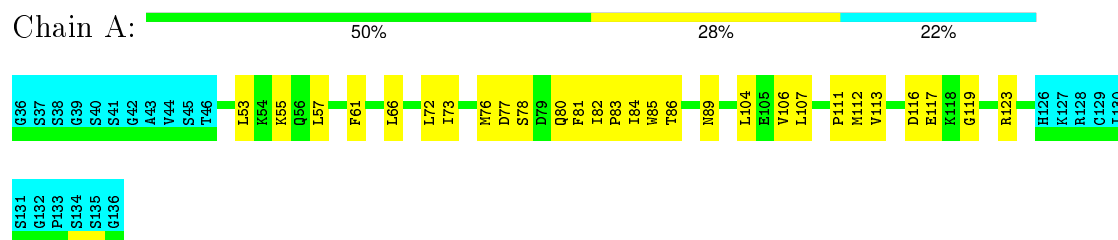
4.2.17 Score per residue for model 17

- Molecule 1: C-Mpl binding protein



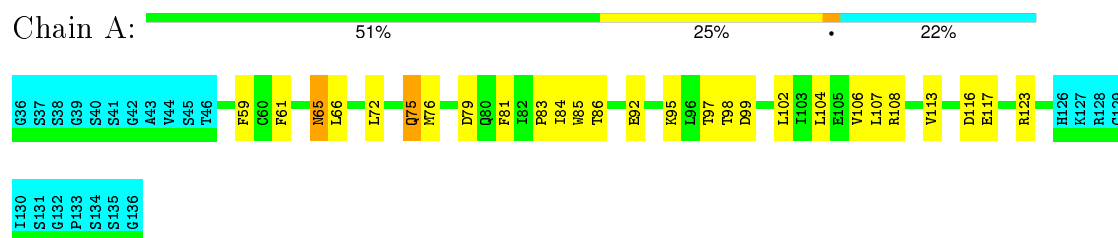
4.2.18 Score per residue for model 18

- Molecule 1: C-Mpl binding protein



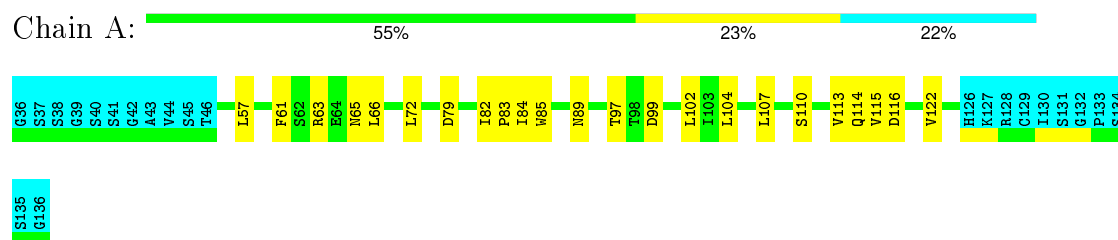
4.2.19 Score per residue for model 19

- Molecule 1: C-Mpl binding protein



4.2.20 Score per residue for model 20

- Molecule 1: C-Mpl binding protein



5 Refinement protocol and experimental data overview ⓘ

The models were refined using the following method: *torsion angle dynamics, simulated annealing*.

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	2.0.17
CYANA	refinement	2.0.17

No chemical shift data was provided. No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality ⓘ

6.1 Standard geometry ⓘ

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	644	654	654	21±4
All	All	12880	13080	13080	423

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 16.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:84:ILE:HD13	1:A:104:LEU:HD21	0.97	1.32	8	17
1:A:53:LEU:HD23	1:A:106:VAL:HG21	0.88	1.46	17	8
1:A:84:ILE:CD1	1:A:104:LEU:HD21	0.83	2.03	8	8
1:A:66:LEU:HD23	1:A:72:LEU:HD23	0.81	1.52	19	20
1:A:107:LEU:O	1:A:113:VAL:HG23	0.79	1.76	13	13
1:A:66:LEU:O	1:A:73:ILE:HD11	0.77	1.80	4	14
1:A:57:LEU:HD13	1:A:107:LEU:HG	0.76	1.57	9	15
1:A:53:LEU:CD2	1:A:106:VAL:HG21	0.76	2.11	17	8
1:A:114:GLN:O	1:A:122:VAL:HG23	0.75	1.81	14	5
1:A:104:LEU:HD22	1:A:115:VAL:HG11	0.74	1.59	20	1
1:A:97:THR:HG21	1:A:102:LEU:HD23	0.74	1.58	8	5
1:A:107:LEU:CD2	1:A:113:VAL:HG21	0.71	2.16	20	5
1:A:84:ILE:HD13	1:A:104:LEU:CD2	0.70	2.17	12	4
1:A:97:THR:HG23	1:A:102:LEU:HD23	0.69	1.62	11	8

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:66:LEU:HD13	1:A:76:MET:HE1	0.68	1.66	2	1
1:A:102:LEU:O	1:A:106:VAL:HG23	0.67	1.90	14	5
1:A:107:LEU:HD23	1:A:113:VAL:HG21	0.66	1.68	20	11
1:A:61:PHE:CE2	1:A:82:ILE:HD12	0.65	2.26	18	4
1:A:97:THR:HG22	1:A:99:ASP:H	0.65	1.52	8	8
1:A:66:LEU:HD22	1:A:76:MET:CE	0.59	2.27	8	3
1:A:57:LEU:HD22	1:A:107:LEU:HD21	0.59	1.73	2	5
1:A:85:TRP:CD2	1:A:86:THR:N	0.59	2.71	19	9
1:A:53:LEU:HD22	1:A:97:THR:OG1	0.57	1.99	11	4
1:A:57:LEU:CD2	1:A:87:VAL:HG11	0.56	2.31	12	3
1:A:76:MET:CE	1:A:82:ILE:HD11	0.56	2.30	3	1
1:A:61:PHE:CD2	1:A:113:VAL:HG11	0.56	2.35	19	1
1:A:115:VAL:HA	1:A:122:VAL:HG22	0.55	1.78	5	1
1:A:59:PHE:CE2	1:A:65:ASN:OD1	0.54	2.61	19	1
1:A:57:LEU:HD12	1:A:106:VAL:CG1	0.54	2.31	8	5
1:A:114:GLN:C	1:A:122:VAL:HG23	0.54	2.23	14	2
1:A:57:LEU:HD21	1:A:87:VAL:HG11	0.54	1.79	12	1
1:A:52:CYS:SG	1:A:96:LEU:HD22	0.53	2.44	6	1
1:A:110:SER:HB2	1:A:113:VAL:HG22	0.52	1.81	11	1
1:A:85:TRP:CZ2	1:A:86:THR:OG1	0.52	2.61	18	3
1:A:85:TRP:CG	1:A:86:THR:N	0.51	2.78	19	14
1:A:82:ILE:HG23	1:A:83:PRO:HD2	0.51	1.83	3	5
1:A:66:LEU:HD23	1:A:72:LEU:CD2	0.51	2.33	8	2
1:A:75:GLN:CD	1:A:75:GLN:C	0.51	2.70	19	1
1:A:66:LEU:CD2	1:A:72:LEU:HD23	0.50	2.30	16	5
1:A:52:CYS:SG	1:A:53:LEU:N	0.50	2.85	6	1
1:A:86:THR:O	1:A:89:ASN:ND2	0.50	2.45	13	2
1:A:101:ASP:OD1	1:A:102:LEU:N	0.50	2.45	7	1
1:A:83:PRO:HB2	1:A:85:TRP:CD1	0.50	2.42	14	20
1:A:107:LEU:O	1:A:113:VAL:CG2	0.50	2.60	15	3
1:A:85:TRP:CE3	1:A:85:TRP:C	0.49	2.85	19	5
1:A:112:MET:O	1:A:125:SER:N	0.49	2.45	14	4
1:A:59:PHE:O	1:A:65:ASN:CB	0.49	2.60	9	10
1:A:85:TRP:C	1:A:85:TRP:CE3	0.49	2.85	18	3
1:A:75:GLN:CD	1:A:76:MET:N	0.49	2.66	19	1
1:A:66:LEU:HD22	1:A:76:MET:HE3	0.49	1.84	8	1
1:A:111:PRO:O	1:A:113:VAL:N	0.49	2.45	12	10
1:A:116:ASP:O	1:A:119:GLY:N	0.49	2.45	10	5
1:A:80:GLN:CG	1:A:80:GLN:O	0.49	2.61	12	1
1:A:105:GLU:O	1:A:109:SER:CB	0.48	2.61	9	4
1:A:81:PHE:CZ	1:A:123:ARG:HD3	0.48	2.44	6	5

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:77:ASP:OD1	1:A:79:ASP:N	0.48	2.46	3	1
1:A:77:ASP:N	1:A:81:PHE:O	0.48	2.47	13	8
1:A:66:LEU:O	1:A:73:ILE:CD1	0.48	2.61	12	5
1:A:85:TRP:CD2	1:A:85:TRP:C	0.48	2.87	19	6
1:A:89:ASN:OD1	1:A:89:ASN:C	0.47	2.53	9	1
1:A:79:ASP:O	1:A:81:PHE:N	0.47	2.47	12	1
1:A:61:PHE:CE1	1:A:82:ILE:HD12	0.47	2.44	6	1
1:A:89:ASN:C	1:A:89:ASN:OD1	0.47	2.53	13	3
1:A:116:ASP:CB	1:A:121:LYS:O	0.47	2.63	5	1
1:A:76:MET:HE3	1:A:82:ILE:HD11	0.47	1.85	3	1
1:A:61:PHE:CZ	1:A:82:ILE:HD12	0.46	2.45	11	1
1:A:85:TRP:C	1:A:85:TRP:CD2	0.46	2.88	6	10
1:A:97:THR:CG2	1:A:102:LEU:HD23	0.46	2.40	3	3
1:A:81:PHE:CD1	1:A:123:ARG:HB3	0.46	2.46	6	7
1:A:51:GLU:O	1:A:55:LYS:CD	0.46	2.63	9	1
1:A:99:ASP:OD1	1:A:101:ASP:CG	0.45	2.55	7	1
1:A:81:PHE:CE2	1:A:123:ARG:HD3	0.45	2.47	6	1
1:A:85:TRP:CE2	1:A:86:THR:OG1	0.45	2.67	7	2
1:A:78:SER:O	1:A:80:GLN:NE2	0.45	2.49	18	1
1:A:94:LYS:O	1:A:98:THR:HG23	0.45	2.11	2	3
1:A:116:ASP:O	1:A:117:GLU:C	0.45	2.55	6	9
1:A:111:PRO:O	1:A:112:MET:C	0.45	2.55	12	5
1:A:81:PHE:CE1	1:A:123:ARG:HB3	0.45	2.46	6	2
1:A:79:ASP:O	1:A:80:GLN:C	0.45	2.54	12	1
1:A:77:ASP:OD1	1:A:77:ASP:C	0.45	2.56	7	3
1:A:113:VAL:C	1:A:114:GLN:OE1	0.45	2.55	14	1
1:A:53:LEU:HD12	1:A:53:LEU:O	0.44	2.12	7	1
1:A:110:SER:HB3	1:A:113:VAL:HG22	0.44	1.90	13	1
1:A:116:ASP:C	1:A:116:ASP:OD1	0.44	2.57	6	1
1:A:114:GLN:O	1:A:122:VAL:HG13	0.44	2.13	5	1
1:A:89:ASN:O	1:A:90:MET:C	0.44	2.56	8	1
1:A:57:LEU:HD12	1:A:106:VAL:HG11	0.43	1.89	8	1
1:A:61:PHE:CD2	1:A:113:VAL:CG1	0.43	3.01	19	1
1:A:85:TRP:O	1:A:89:ASN:ND2	0.43	2.51	14	1
1:A:81:PHE:CE2	1:A:123:ARG:HD2	0.43	2.48	7	1
1:A:85:TRP:CE3	1:A:86:THR:HA	0.43	2.49	18	1
1:A:114:GLN:OE1	1:A:115:VAL:C	0.42	2.57	20	1
1:A:81:PHE:CZ	1:A:123:ARG:NH2	0.42	2.86	8	1
1:A:104:LEU:HD22	1:A:115:VAL:CG1	0.42	2.36	20	1
1:A:107:LEU:HD22	1:A:113:VAL:HG21	0.42	1.91	20	1
1:A:68:LYS:CD	1:A:68:LYS:C	0.42	2.87	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:111:PRO:C	1:A:113:VAL:N	0.42	2.72	18	6
1:A:89:ASN:O	1:A:94:LYS:CD	0.42	2.68	5	1
1:A:60:CYS:SG	1:A:90:MET:CE	0.42	3.07	16	1
1:A:110:SER:O	1:A:113:VAL:HG22	0.42	2.14	17	1
1:A:66:LEU:O	1:A:73:ILE:CG1	0.42	2.68	14	1
1:A:81:PHE:CZ	1:A:123:ARG:CD	0.42	3.02	7	1
1:A:57:LEU:HD13	1:A:107:LEU:CD2	0.41	2.45	18	1
1:A:63:ARG:CZ	1:A:63:ARG:HB2	0.41	2.46	16	1
1:A:51:GLU:HG3	1:A:52:CYS:N	0.41	2.30	9	1
1:A:55:LYS:HG3	1:A:56:GLN:N	0.41	2.30	10	1
1:A:115:VAL:O	1:A:116:ASP:C	0.41	2.59	20	1
1:A:110:SER:CB	1:A:113:VAL:HG22	0.41	2.45	11	1
1:A:90:MET:HB2	1:A:93:ILE:CG1	0.41	2.46	5	2
1:A:48:ASP:OD1	1:A:48:ASP:C	0.40	2.60	16	1
1:A:53:LEU:HD11	1:A:57:LEU:HD11	0.40	1.93	9	2
1:A:110:SER:O	1:A:113:VAL:O	0.40	2.39	20	2
1:A:56:GLN:OE1	1:A:92:GLU:OE1	0.40	2.40	4	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	79/101 (78%)	73±2 (92±3%)	5±2 (6±2%)	1±1 (2±1%)	17	61
All	All	1580/2020 (78%)	1460 (92%)	94 (6%)	26 (2%)	17	61

All 5 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	76	MET	13
1	A	112	MET	9
1	A	125	SER	2
1	A	80	GLN	1
1	A	116	ASP	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	77/93 (83%)	74±1 (96±2%)	3±1 (4±2%)	39 82
All	All	1540/1860 (83%)	1473 (96%)	67 (4%)	39 82

All 19 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	65	ASN	12
1	A	89	ASN	11
1	A	77	ASP	7
1	A	79	ASP	6
1	A	50	LYS	5
1	A	95	LYS	4
1	A	98	THR	4
1	A	118	LYS	2
1	A	68	LYS	2
1	A	108	ARG	2
1	A	92	GLU	2
1	A	55	LYS	2
1	A	117	GLU	2
1	A	48	ASP	1
1	A	101	ASP	1
1	A	91	GLU	1
1	A	63	ARG	1
1	A	80	GLN	1
1	A	75	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 [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

No chemical shift data were provided