



Full wwPDB NMR Structure Validation Report ⓘ

Feb 13, 2017 – 12:22 am GMT

PDB ID : 2LS8
Title : Solution structure of human C-type lectin domain family 4 member D
Authors : Harris, R.; Gaudette, J.; Bandaranayake, A.D.; Banu, R.; Bonanno, J.B.; Calarese, D.A.; Celikgil, A.; Chamala, S.; Chan, M.K.; Chaparro, R.; Evans, B.; Garforth, S.; Gizzi, A.; Hillerich, B.; Kar, A.; Lafleur, J.; Lim, S.; Love, J.; Matikainen, B.; Patel, H.; Seidel, R.D.; Smith, B.; Stead, M.; Girvin, M.E.; Almo, S.C.; New York Structural Genomics Research Consortium (NYSGRG)
Deposited on : 2012-04-23

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

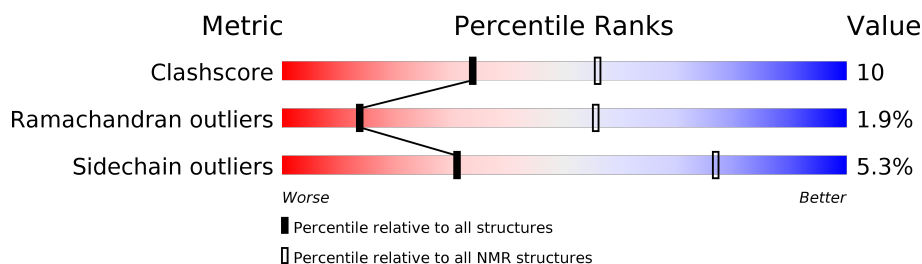
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 90%.

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	156	

2 Ensemble composition and analysis

This entry contains 20 models. Model 18 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:3-A:89, A:100-A:130 (118)	0.53	18
2	A:131-A:142 (12)	0.35	9

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 2 single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2273 atoms, of which 1096 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called C-type lectin domain family 4 member D.

Mol	Chain	Residues	Atoms						Trace
1	A	142	Total	C	H	N	O	S	0
			2273	744	1096	207	217	9	

There are 24 discrepancies between the modelled and reference sequences:

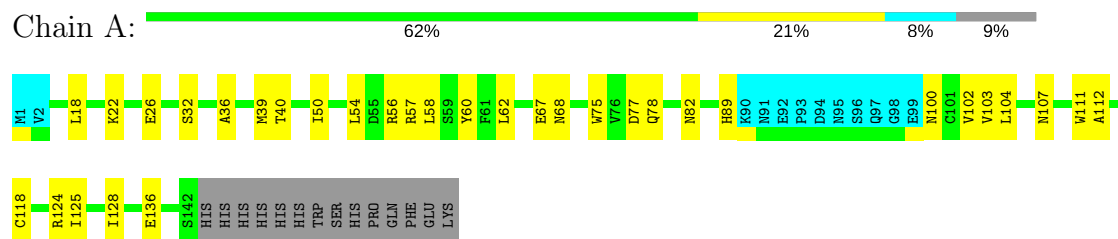
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP Q8WXI8
A	2	VAL	-	EXPRESSION TAG	UNP Q8WXI8
A	135	ALA	-	EXPRESSION TAG	UNP Q8WXI8
A	136	GLU	-	EXPRESSION TAG	UNP Q8WXI8
A	137	ASN	-	EXPRESSION TAG	UNP Q8WXI8
A	138	LEU	-	EXPRESSION TAG	UNP Q8WXI8
A	139	TYR	-	EXPRESSION TAG	UNP Q8WXI8
A	140	PHE	-	EXPRESSION TAG	UNP Q8WXI8
A	141	GLN	-	EXPRESSION TAG	UNP Q8WXI8
A	142	SER	-	EXPRESSION TAG	UNP Q8WXI8
A	143	HIS	-	EXPRESSION TAG	UNP Q8WXI8
A	144	HIS	-	EXPRESSION TAG	UNP Q8WXI8
A	145	HIS	-	EXPRESSION TAG	UNP Q8WXI8
A	146	HIS	-	EXPRESSION TAG	UNP Q8WXI8
A	147	HIS	-	EXPRESSION TAG	UNP Q8WXI8
A	148	HIS	-	EXPRESSION TAG	UNP Q8WXI8
A	149	TRP	-	EXPRESSION TAG	UNP Q8WXI8
A	150	SER	-	EXPRESSION TAG	UNP Q8WXI8
A	151	HIS	-	EXPRESSION TAG	UNP Q8WXI8
A	152	PRO	-	EXPRESSION TAG	UNP Q8WXI8
A	153	GLN	-	EXPRESSION TAG	UNP Q8WXI8
A	154	PHE	-	EXPRESSION TAG	UNP Q8WXI8
A	155	GLU	-	EXPRESSION TAG	UNP Q8WXI8
A	156	LYS	-	EXPRESSION TAG	UNP Q8WXI8

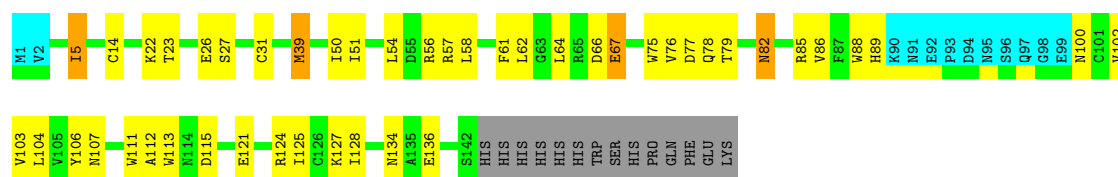
4 Residue-property plots [i](#)

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: C-type lectin domain family 4 member D

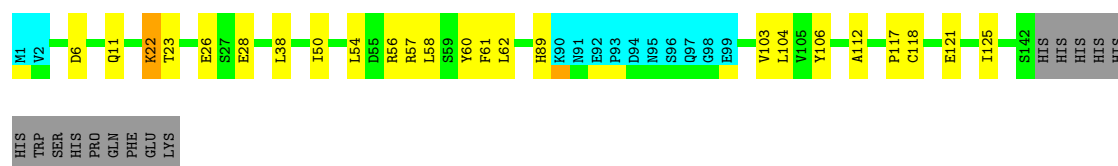




4.2.3 Score per residue for model 3

- Molecule 1: C-type lectin domain family 4 member D

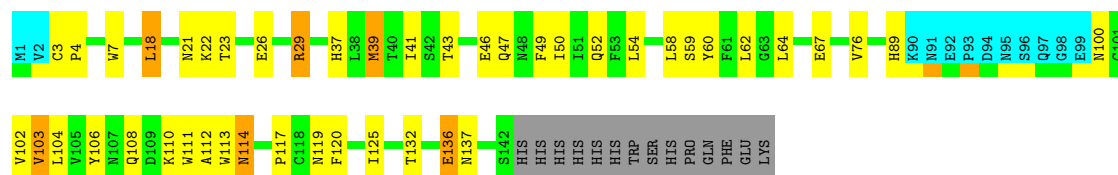
Chain A: 68% 15% 8% 9%



4.2.4 Score per residue for model 4

- Molecule 1: C-type lectin domain family 4 member D

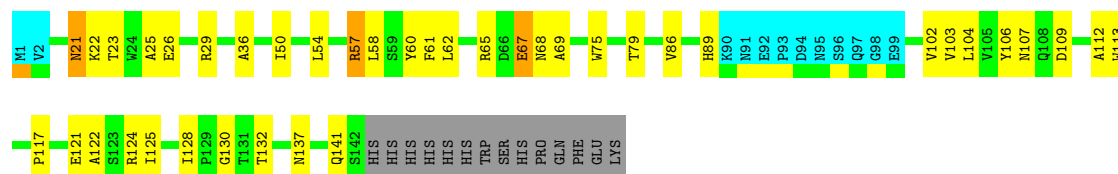
Chain A: 54% 25% 8% 9%



4.2.5 Score per residue for model 5

- Molecule 1: C-type lectin domain family 4 member D

Chain A: 58% 24% 8% 9%



4.2.6 Score per residue for model 6

- Molecule 1: C-type lectin domain family 4 member D

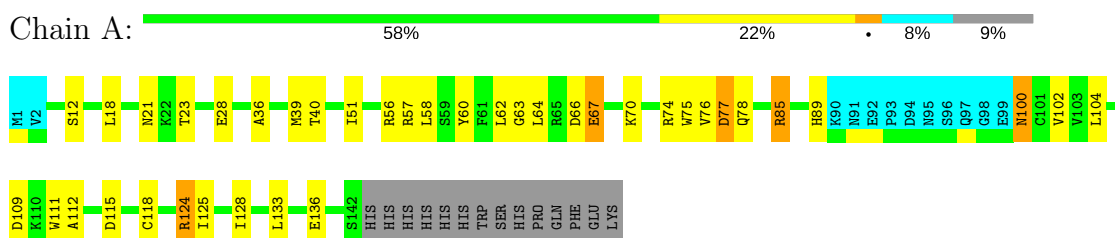
- Molecule 1: C-type lectin domain family 4 member D

4.2.8 Score per residue for model 8

- Molecule 1: C-type lectin domain family 4 member D

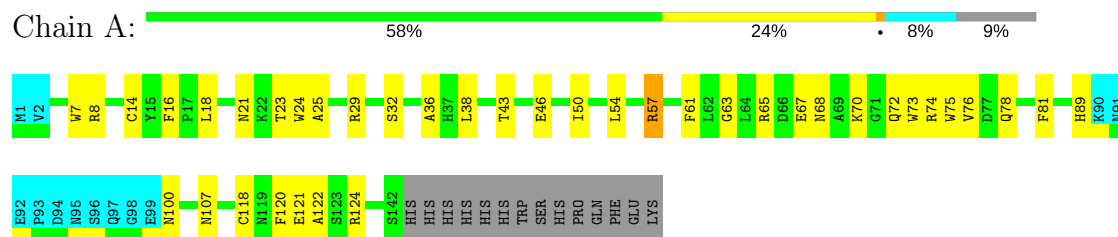
4.2.9 Score per residue for model 9

- Molecule 1: C-type lectin domain family 4 member D



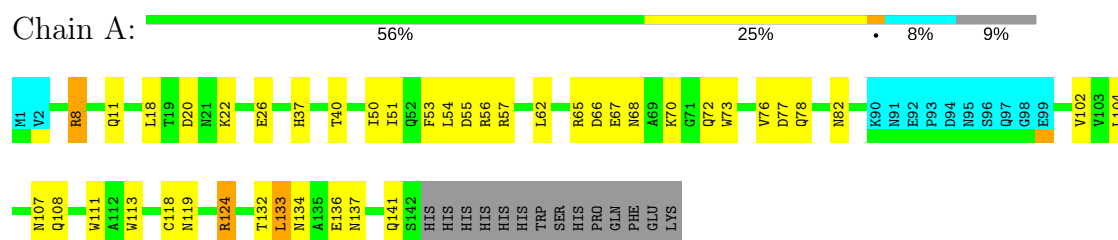
4.2.10 Score per residue for model 10

- Molecule 1: C-type lectin domain family 4 member D



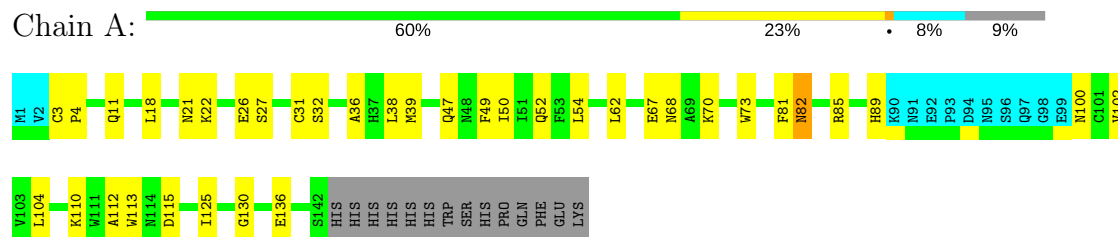
4.2.11 Score per residue for model 11

- Molecule 1: C-type lectin domain family 4 member D



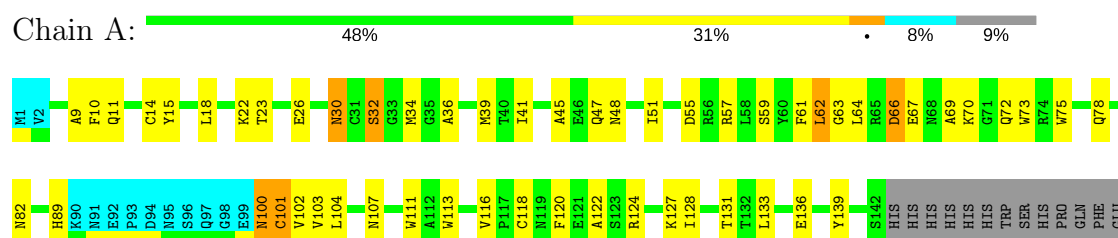
4.2.12 Score per residue for model 12

- Molecule 1: C-type lectin domain family 4 member D



4.2.13 Score per residue for model 13

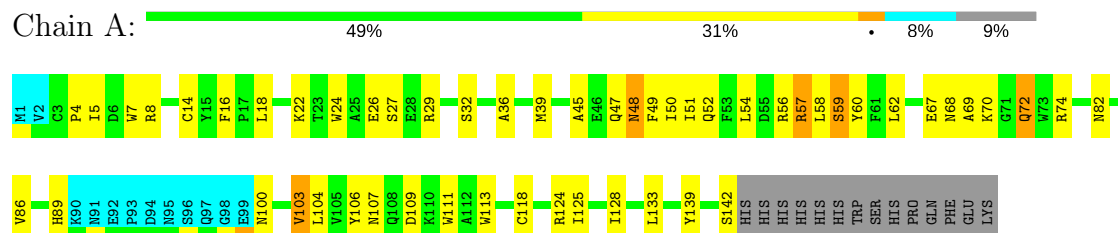
- Molecule 1: C-type lectin domain family 4 member D



LYS

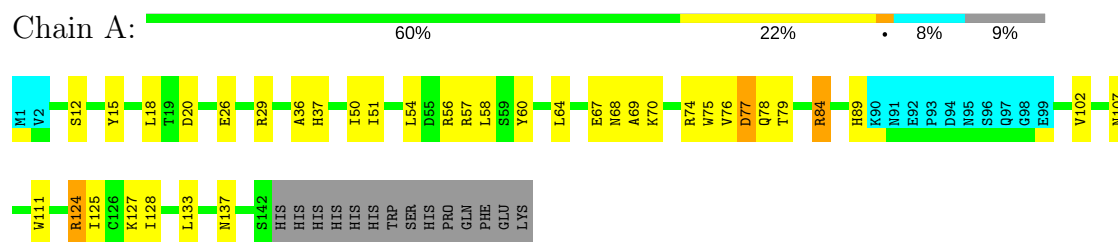
4.2.14 Score per residue for model 14

- Molecule 1: C-type lectin domain family 4 member D



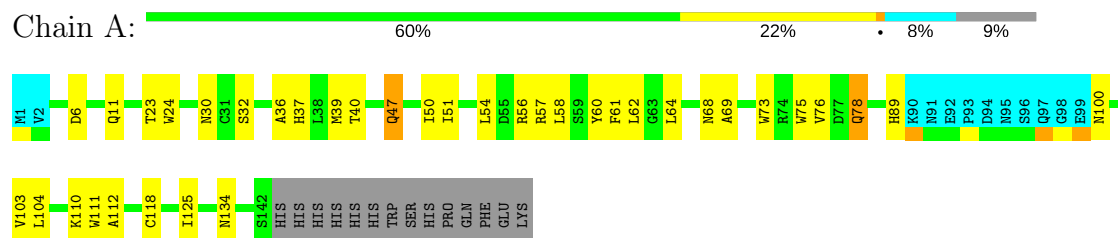
4.2.15 Score per residue for model 15

- Molecule 1: C-type lectin domain family 4 member D



4.2.16 Score per residue for model 16

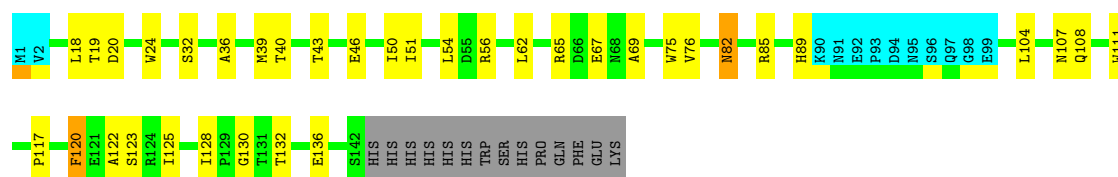
- Molecule 1: C-type lectin domain family 4 member D



4.2.17 Score per residue for model 17

- Molecule 1: C-type lectin domain family 4 member D

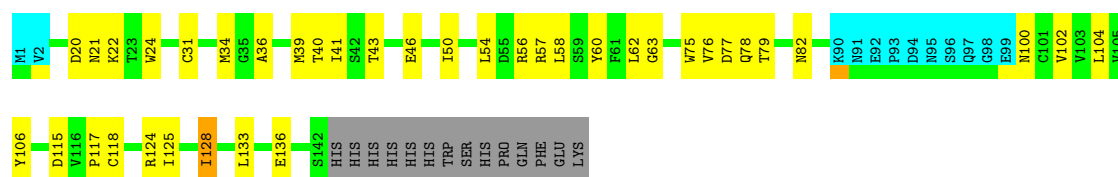




4.2.18 Score per residue for model 18 (medoid)

- Molecule 1: C-type lectin domain family 4 member D

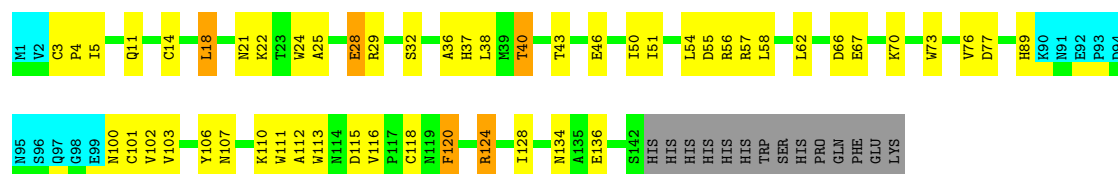
Chain A: 59% 24% 8% 9%



4.2.19 Score per residue for model 19

- Molecule 1: C-type lectin domain family 4 member D

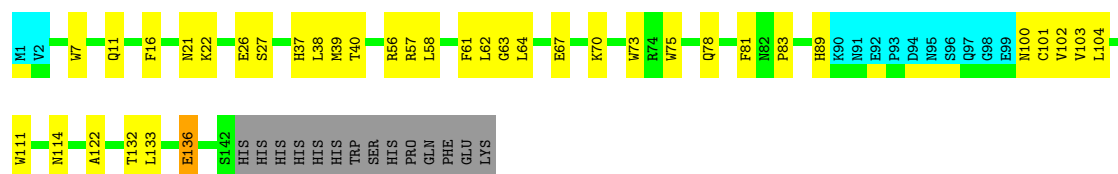
Chain A: 50% 30% 8% 9%



4.2.20 Score per residue for model 20

- Molecule 1: C-type lectin domain family 4 member D

Chain A: 60% 23% 8% 9%



5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *20 structures for lowest energy*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
CNS	structure solution	1.21
CNS	refinement	1.21
ARIA	structure solution	2.3

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)	2ls8_cs.str
Number of chemical shift lists	1
Total number of shifts	1889
Number of shifts mapped to atoms	1889
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	90%

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	1085	1012	1010	22±5
All	All	21700	20240	20200	431

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 10.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:22:LYS:HD3	1:A:26:GLU:HB3	0.85	1.48	3	2
1:A:67:GLU:HG2	1:A:70:LYS:HD2	0.79	1.51	9	1
1:A:39:MET:HB2	1:A:125:ILE:HB	0.77	1.54	1	8
1:A:45:ALA:HA	1:A:48:ASN:ND2	0.74	1.97	1	3
1:A:64:LEU:HB2	1:A:102:VAL:HG12	0.72	1.58	2	1
1:A:32:SER:HA	1:A:36:ALA:O	0.72	1.82	12	9
1:A:57:ARG:HG2	1:A:58:LEU:HD22	0.71	1.61	3	10
1:A:22:LYS:HB2	1:A:124:ARG:HD2	0.70	1.62	18	2
1:A:59:SER:HB3	1:A:103:VAL:HG11	0.70	1.63	4	1
1:A:133:LEU:H	1:A:133:LEU:HD13	0.69	1.45	11	1
1:A:51:ILE:HB	1:A:111:TRP:CD1	0.69	2.22	9	8
1:A:100:ASN:HB2	1:A:115:ASP:OD1	0.66	1.91	12	2
1:A:3:CYS:SG	1:A:4:PRO:HD2	0.66	2.30	19	1
1:A:22:LYS:HB2	1:A:124:ARG:HD3	0.64	1.70	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:37:HIS:CD2	1:A:77:ASP:HA	0.64	2.27	1	1
1:A:102:VAL:HB	1:A:113:TRP:HB3	0.64	1.69	19	6
1:A:56:ARG:HA	1:A:111:TRP:CZ2	0.63	2.28	19	8
1:A:36:ALA:HB2	1:A:128:ILE:HG12	0.63	1.70	18	1
1:A:62:LEU:HD11	1:A:104:LEU:HG	0.62	1.69	20	5
1:A:21:ASN:HB3	1:A:121:GLU:HB3	0.62	1.69	8	2
1:A:22:LYS:HD2	1:A:26:GLU:HB3	0.62	1.70	20	3
1:A:59:SER:HB3	1:A:103:VAL:HG21	0.61	1.71	13	2
1:A:61:PHE:HA	1:A:103:VAL:HG12	0.61	1.70	3	8
1:A:47:GLN:HE22	1:A:104:LEU:HD11	0.61	1.54	12	1
1:A:66:ASP:HB3	1:A:100:ASN:HB3	0.60	1.73	6	1
1:A:56:ARG:HG3	1:A:106:TYR:CD2	0.60	2.31	19	1
1:A:18:LEU:HB2	1:A:124:ARG:HG3	0.60	1.74	11	4
1:A:40:THR:HB	1:A:76:VAL:HG12	0.59	1.72	19	1
1:A:101:CYS:HB2	1:A:116:VAL:O	0.59	1.98	13	1
1:A:62:LEU:HD11	1:A:104:LEU:HB2	0.59	1.75	4	1
1:A:21:ASN:O	1:A:22:LYS:HD2	0.59	1.97	18	4
1:A:4:PRO:HB2	1:A:7:TRP:CD1	0.59	2.33	8	3
1:A:67:GLU:O	1:A:70:LYS:HG3	0.58	1.98	20	4
1:A:103:VAL:O	1:A:113:TRP:HA	0.58	1.98	4	2
1:A:36:ALA:HB2	1:A:128:ILE:HB	0.58	1.76	17	4
1:A:47:GLN:HE22	1:A:51:ILE:HD13	0.58	1.59	16	1
1:A:133:LEU:HD23	1:A:133:LEU:H	0.58	1.58	20	1
1:A:73:TRP:CH2	1:A:100:ASN:HB3	0.58	2.33	13	1
1:A:62:LEU:HD21	1:A:104:LEU:HG	0.57	1.75	1	5
1:A:62:LEU:HB2	1:A:102:VAL:HG13	0.57	1.76	9	2
1:A:15:TYR:HE2	1:A:39:MET:HG2	0.57	1.60	8	1
1:A:66:ASP:HB2	1:A:101:CYS:HA	0.57	1.76	19	1
1:A:89:HIS:CE1	1:A:107:ASN:HB3	0.56	2.35	6	7
1:A:8:ARG:HD3	1:A:53:PHE:CE1	0.56	2.35	11	1
1:A:67:GLU:CD	1:A:67:GLU:H	0.56	2.04	2	1
1:A:139:TYR:HA	1:A:142:SER:OG	0.56	2.00	1	2
1:A:23:THR:HG23	1:A:121:GLU:OE2	0.56	2.01	5	2
1:A:100:ASN:HB2	1:A:115:ASP:OD2	0.56	2.01	19	1
1:A:136:GLU:CD	1:A:137:ASN:H	0.56	2.04	8	3
1:A:24:TRP:CH2	1:A:65:ARG:HG3	0.56	2.36	17	1
1:A:60:TYR:HB3	1:A:125:ILE:HG13	0.55	1.78	5	10
1:A:132:THR:O	1:A:136:GLU:HG2	0.55	2.02	11	1
1:A:14:CYS:HB2	1:A:128:ILE:HB	0.55	1.79	19	2
1:A:64:LEU:HD23	1:A:64:LEU:H	0.55	1.62	8	1
1:A:84:ARG:NE	1:A:84:ARG:HA	0.55	2.16	15	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:49:PHE:O	1:A:52:GLN:HG2	0.55	2.02	7	4
1:A:64:LEU:HB2	1:A:102:VAL:HG22	0.54	1.78	4	1
1:A:37:HIS:CD2	1:A:77:ASP:HB2	0.54	2.37	19	1
1:A:75:TRP:HB2	1:A:79:THR:HB	0.54	1.80	2	4
1:A:89:HIS:CE1	1:A:107:ASN:HB2	0.54	2.38	14	2
1:A:62:LEU:CD2	1:A:104:LEU:HG	0.54	2.32	13	1
1:A:63:GLY:O	1:A:75:TRP:HA	0.54	2.01	18	4
1:A:66:ASP:HA	1:A:70:LYS:HE2	0.54	1.79	8	1
1:A:114:ASN:HD22	1:A:114:ASN:C	0.54	2.07	4	1
1:A:24:TRP:HB3	1:A:118:CYS:HB3	0.54	1.80	16	2
1:A:73:TRP:O	1:A:81:PHE:HB2	0.53	2.03	20	2
1:A:24:TRP:CH2	1:A:65:ARG:HB2	0.53	2.39	10	2
1:A:25:ALA:O	1:A:29:ARG:HG2	0.53	2.03	8	3
1:A:82:ASN:ND2	1:A:85:ARG:HB2	0.53	2.18	2	1
1:A:66:ASP:HB3	1:A:100:ASN:HB2	0.53	1.79	13	1
1:A:23:THR:HA	1:A:120:PHE:O	0.53	2.04	13	3
1:A:89:HIS:CE1	1:A:112:ALA:HB3	0.52	2.39	2	5
1:A:37:HIS:HB3	1:A:76:VAL:HG13	0.52	1.80	16	1
1:A:73:TRP:CH2	1:A:102:VAL:HG13	0.52	2.38	13	2
1:A:89:HIS:CD2	1:A:112:ALA:HB3	0.52	2.40	12	5
1:A:22:LYS:HE2	1:A:26:GLU:HB3	0.52	1.82	5	2
1:A:84:ARG:HE	1:A:84:ARG:HA	0.52	1.65	15	1
1:A:47:GLN:HE21	1:A:104:LEU:HD21	0.51	1.64	13	1
1:A:55:ASP:OD1	1:A:57:ARG:HB3	0.51	2.06	11	3
1:A:23:THR:HB	1:A:118:CYS:O	0.51	2.06	16	3
1:A:133:LEU:H	1:A:133:LEU:HD12	0.51	1.65	15	1
1:A:106:TYR:CE1	1:A:109:ASP:HA	0.51	2.40	5	1
1:A:22:LYS:HD3	1:A:26:GLU:CB	0.51	2.33	12	1
1:A:43:THR:O	1:A:46:GLU:HB3	0.51	2.05	6	6
1:A:40:THR:HA	1:A:75:TRP:CE3	0.51	2.41	18	4
1:A:74:ARG:HD3	1:A:78:GLN:HA	0.51	1.82	15	1
1:A:50:ILE:O	1:A:54:LEU:HG	0.50	2.05	1	17
1:A:124:ARG:HH11	1:A:124:ARG:HB2	0.50	1.65	19	1
1:A:64:LEU:HB2	1:A:102:VAL:CG1	0.50	2.36	9	1
1:A:85:ARG:N	1:A:85:ARG:HD2	0.50	2.21	9	1
1:A:40:THR:HG21	1:A:77:ASP:HB3	0.50	1.83	18	1
1:A:89:HIS:HA	1:A:112:ALA:CB	0.50	2.37	8	3
1:A:74:ARG:NH2	1:A:78:GLN:HA	0.50	2.20	10	1
1:A:70:LYS:HB3	1:A:72:GLN:OE1	0.50	2.07	14	1
1:A:22:LYS:HA	1:A:121:GLU:OE1	0.50	2.06	5	1
1:A:22:LYS:HD2	1:A:26:GLU:CB	0.50	2.37	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:10:PHE:HB3	1:A:15:TYR:CE1	0.50	2.42	6	2
1:A:132:THR:HB	1:A:136:GLU:HG2	0.49	1.84	11	1
1:A:66:ASP:HA	1:A:70:LYS:HD3	0.49	1.84	9	1
1:A:67:GLU:HB3	1:A:70:LYS:HZ3	0.49	1.67	10	1
1:A:39:MET:SD	1:A:125:ILE:HG21	0.49	2.48	12	1
1:A:18:LEU:HD23	1:A:18:LEU:N	0.49	2.22	19	2
1:A:9:ALA:HB1	1:A:133:LEU:HD21	0.49	1.83	13	1
1:A:64:LEU:HD21	1:A:102:VAL:HG22	0.49	1.83	6	1
1:A:18:LEU:HG	1:A:124:ARG:NH2	0.48	2.23	13	1
1:A:62:LEU:HD13	1:A:64:LEU:HD23	0.48	1.85	13	1
1:A:25:ALA:O	1:A:29:ARG:HG3	0.48	2.07	10	2
1:A:3:CYS:HB3	1:A:4:PRO:HD2	0.48	1.83	4	2
1:A:28:GLU:HG2	1:A:38:LEU:HD13	0.48	1.86	3	1
1:A:56:ARG:HD3	1:A:106:TYR:CE1	0.47	2.43	14	5
1:A:20:ASP:OD2	1:A:22:LYS:HD3	0.47	2.09	18	1
1:A:108:GLN:O	1:A:110:LYS:HG2	0.47	2.10	4	1
1:A:31:CYS:HA	1:A:34:MET:SD	0.47	2.49	8	2
1:A:28:GLU:HG2	1:A:38:LEU:HB2	0.47	1.86	19	1
1:A:3:CYS:HB2	1:A:7:TRP:HB2	0.47	1.86	6	1
1:A:66:ASP:HB3	1:A:100:ASN:O	0.47	2.10	9	1
1:A:27:SER:HB3	1:A:124:ARG:HD3	0.47	1.87	14	1
1:A:89:HIS:NE2	1:A:107:ASN:HB3	0.47	2.25	19	1
1:A:8:ARG:CZ	1:A:8:ARG:HA	0.47	2.40	11	1
1:A:106:TYR:CZ	1:A:109:ASP:HA	0.47	2.45	5	1
1:A:40:THR:HG22	1:A:75:TRP:HB3	0.47	1.86	9	2
1:A:61:PHE:CE2	1:A:122:ALA:HB3	0.47	2.45	20	3
1:A:18:LEU:N	1:A:18:LEU:HD23	0.47	2.25	4	1
1:A:39:MET:CE	1:A:41:ILE:HD13	0.46	2.40	13	2
1:A:82:ASN:HB2	1:A:85:ARG:HB2	0.46	1.86	17	1
1:A:39:MET:HA	1:A:127:LYS:HB3	0.46	1.87	13	1
1:A:37:HIS:CD2	1:A:76:VAL:HG13	0.46	2.46	11	2
1:A:59:SER:HB3	1:A:103:VAL:CG1	0.46	2.38	4	1
1:A:18:LEU:HD22	1:A:124:ARG:NH2	0.46	2.25	14	1
1:A:67:GLU:CB	1:A:70:LYS:HD2	0.46	2.41	14	1
1:A:137:ASN:O	1:A:141:GLN:HG2	0.46	2.10	5	3
1:A:38:LEU:HB2	1:A:76:VAL:HG21	0.46	1.88	10	1
1:A:40:THR:HG21	1:A:77:ASP:OD1	0.46	2.11	11	1
1:A:18:LEU:HB3	1:A:20:ASP:OD1	0.46	2.10	15	1
1:A:64:LEU:HB3	1:A:102:VAL:CG2	0.45	2.42	15	1
1:A:82:ASN:CB	1:A:85:ARG:HB2	0.45	2.42	8	3
1:A:14:CYS:HB2	1:A:128:ILE:HG23	0.45	1.88	13	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:41:ILE:HA	1:A:46:GLU:OE2	0.45	2.11	18	1
1:A:67:GLU:O	1:A:67:GLU:HG3	0.45	2.10	17	1
1:A:67:GLU:HB3	1:A:70:LYS:HD2	0.45	1.88	14	1
1:A:77:ASP:O	1:A:78:GLN:HB3	0.45	2.11	8	1
1:A:39:MET:CB	1:A:125:ILE:HB	0.45	2.40	6	1
1:A:27:SER:HA	1:A:124:ARG:HD2	0.45	1.87	2	1
1:A:24:TRP:CZ2	1:A:76:VAL:HG23	0.45	2.47	19	1
1:A:26:GLU:HG2	1:A:29:ARG:HH11	0.45	1.71	15	1
1:A:28:GLU:OE2	1:A:76:VAL:HG11	0.45	2.12	9	1
1:A:133:LEU:HD22	1:A:134:ASN:H	0.45	1.71	11	1
1:A:26:GLU:O	1:A:30:ASN:HB2	0.45	2.11	13	1
1:A:133:LEU:HD12	1:A:133:LEU:H	0.45	1.72	18	1
1:A:135:ALA:HB1	1:A:138:LEU:HB3	0.44	1.89	6	1
1:A:14:CYS:HB2	1:A:128:ILE:CG2	0.44	2.42	7	2
1:A:66:ASP:HA	1:A:73:TRP:HZ3	0.44	1.72	19	1
1:A:8:ARG:O	1:A:14:CYS:HA	0.44	2.13	14	4
1:A:40:THR:CG2	1:A:75:TRP:HB3	0.44	2.42	16	2
1:A:66:ASP:HB3	1:A:100:ASN:OD1	0.44	2.13	2	1
1:A:104:LEU:HG	1:A:111:TRP:CE3	0.44	2.48	17	1
1:A:8:ARG:HD3	1:A:53:PHE:CD1	0.44	2.48	11	1
1:A:67:GLU:HB3	1:A:70:LYS:NZ	0.44	2.27	10	1
1:A:89:HIS:HA	1:A:112:ALA:HB1	0.44	1.88	7	1
1:A:22:LYS:HZ3	1:A:22:LYS:HB2	0.44	1.72	7	1
1:A:64:LEU:HB2	1:A:102:VAL:HG11	0.44	1.88	9	1
1:A:40:THR:HA	1:A:75:TRP:HZ3	0.44	1.73	8	1
1:A:86:VAL:HG23	1:A:113:TRP:CE3	0.44	2.47	14	1
1:A:7:TRP:CD2	1:A:16:PHE:HB2	0.44	2.48	20	3
1:A:64:LEU:HB3	1:A:102:VAL:HG22	0.44	1.88	20	2
1:A:70:LYS:HE2	1:A:72:GLN:HB2	0.43	1.89	11	1
1:A:31:CYS:O	1:A:32:SER:HB2	0.43	2.13	8	2
1:A:132:THR:O	1:A:136:GLU:HB3	0.43	2.13	20	1
1:A:39:MET:HA	1:A:127:LYS:HB2	0.43	1.90	6	1
1:A:22:LYS:HA	1:A:121:GLU:OE2	0.43	2.12	3	1
1:A:18:LEU:O	1:A:123:SER:HB3	0.43	2.12	17	1
1:A:47:GLN:O	1:A:51:ILE:HG23	0.43	2.13	13	1
1:A:27:SER:CB	1:A:124:ARG:HD3	0.43	2.43	14	1
1:A:63:GLY:CA	1:A:76:VAL:HG23	0.43	2.43	9	1
1:A:62:LEU:HD23	1:A:102:VAL:HG12	0.43	1.89	18	1
1:A:62:LEU:HD21	1:A:104:LEU:N	0.43	2.28	18	1
1:A:64:LEU:HA	1:A:75:TRP:HA	0.43	1.91	6	1
1:A:26:GLU:HG2	1:A:29:ARG:NH1	0.43	2.29	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:22:LYS:HG3	1:A:26:GLU:HB2	0.43	1.90	7	1
1:A:18:LEU:HG	1:A:124:ARG:CZ	0.43	2.44	13	1
1:A:38:LEU:HB3	1:A:76:VAL:HG21	0.43	1.91	19	1
1:A:21:ASN:O	1:A:22:LYS:HG2	0.43	2.14	12	1
1:A:74:ARG:HG2	1:A:75:TRP:N	0.43	2.29	9	1
1:A:103:VAL:HG11	1:A:120:PHE:HE1	0.43	1.73	19	1
1:A:28:GLU:HA	1:A:28:GLU:OE1	0.43	2.14	7	1
1:A:81:PHE:O	1:A:83:PRO:HD3	0.43	2.13	20	2
1:A:89:HIS:HA	1:A:112:ALA:HB3	0.43	1.90	8	2
1:A:86:VAL:HG13	1:A:88:TRP:CD1	0.42	2.49	2	1
1:A:75:TRP:O	1:A:77:ASP:N	0.42	2.53	2	2
1:A:59:SER:CB	1:A:103:VAL:HG21	0.42	2.44	13	1
1:A:64:LEU:HG	1:A:73:TRP:HE3	0.42	1.74	16	1
1:A:66:ASP:HB2	1:A:101:CYS:CA	0.42	2.45	19	1
1:A:5:ILE:H	1:A:5:ILE:HD13	0.42	1.74	2	1
1:A:65:ARG:HG3	1:A:101:CYS:SG	0.42	2.54	7	1
1:A:63:GLY:HA2	1:A:76:VAL:HG23	0.42	1.91	9	1
1:A:106:TYR:HD2	1:A:111:TRP:NE1	0.42	2.12	4	1
1:A:39:MET:SD	1:A:41:ILE:HD13	0.42	2.55	8	1
1:A:66:ASP:CB	1:A:100:ASN:HB2	0.42	2.43	13	1
1:A:36:ALA:HB2	1:A:128:ILE:CG1	0.42	2.44	14	1
1:A:24:TRP:HZ3	1:A:65:ARG:HE	0.42	1.57	10	1
1:A:73:TRP:HB2	1:A:81:PHE:HB2	0.42	1.92	10	2
1:A:40:THR:HG22	1:A:75:TRP:CE3	0.42	2.50	17	1
1:A:64:LEU:HG	1:A:75:TRP:CE2	0.42	2.50	1	1
1:A:43:THR:OG1	1:A:46:GLU:HB2	0.42	2.15	18	1
1:A:24:TRP:HB3	1:A:118:CYS:SG	0.42	2.55	14	1
1:A:39:MET:HA	1:A:127:LYS:CB	0.42	2.45	6	1
1:A:15:TYR:CE2	1:A:127:LYS:HB3	0.42	2.50	7	1
1:A:124:ARG:NH1	1:A:124:ARG:HB2	0.42	2.30	19	1
1:A:65:ARG:HD3	1:A:67:GLU:OE1	0.41	2.15	5	1
1:A:27:SER:OG	1:A:38:LEU:HD13	0.41	2.14	12	1
1:A:62:LEU:HD21	1:A:104:LEU:HB2	0.41	1.91	17	1
1:A:21:ASN:HA	1:A:122:ALA:O	0.41	2.15	5	1
1:A:60:TYR:N	1:A:103:VAL:HG22	0.41	2.30	4	1
1:A:22:LYS:NZ	1:A:22:LYS:HA	0.41	2.30	12	1
1:A:120:PHE:CZ	1:A:122:ALA:HB2	0.41	2.51	17	1
1:A:67:GLU:O	1:A:68:ASN:HB2	0.41	2.16	14	1
1:A:32:SER:N	1:A:36:ALA:O	0.41	2.52	8	1
1:A:65:ARG:HG2	1:A:66:ASP:N	0.41	2.29	11	1
1:A:57:ARG:HG2	1:A:58:LEU:HD13	0.41	1.92	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:26:GLU:HA	1:A:29:ARG:HB2	0.41	1.92	4	1
1:A:83:PRO:O	1:A:84:ARG:HD2	0.41	2.16	7	1
1:A:27:SER:HB3	1:A:38:LEU:HD21	0.41	1.93	20	1
1:A:107:ASN:O	1:A:108:GLN:HB2	0.41	2.16	11	1
1:A:86:VAL:HG23	1:A:113:TRP:CZ3	0.41	2.51	5	1
1:A:113:TRP:CD1	1:A:113:TRP:N	0.41	2.87	8	2
1:A:102:VAL:HG21	1:A:113:TRP:HB3	0.41	1.92	2	1
1:A:15:TYR:CE2	1:A:127:LYS:HE3	0.41	2.51	15	1
1:A:70:LYS:HB2	1:A:72:GLN:HG2	0.41	1.91	13	1
1:A:4:PRO:HD2	1:A:7:TRP:CD1	0.41	2.50	14	1
1:A:57:ARG:HB3	1:A:58:LEU:HD22	0.41	1.91	9	1
1:A:55:ASP:OD2	1:A:57:ARG:HB3	0.41	2.15	19	1
1:A:67:GLU:HG2	1:A:68:ASN:HD22	0.40	1.75	10	1
1:A:56:ARG:HD3	1:A:106:TYR:CZ	0.40	2.51	7	1
1:A:18:LEU:HD22	1:A:124:ARG:NH1	0.40	2.30	15	1
1:A:133:LEU:HD22	1:A:133:LEU:H	0.40	1.75	14	1
1:A:135:ALA:HA	1:A:138:LEU:HB3	0.40	1.92	1	1
1:A:133:LEU:N	1:A:133:LEU:HD12	0.40	2.32	9	1
1:A:23:THR:OG1	1:A:26:GLU:HG2	0.40	2.16	7	1
1:A:5:ILE:HD12	1:A:5:ILE:H	0.40	1.75	7	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	129/156 (83%)	114±3 (88±3%)	13±3 (10±2%)	2±1 (2±1%)	14	55
All	All	2580/3120 (83%)	2278 (88%)	253 (10%)	49 (2%)	14	55

All 19 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	11	GLN	9
1	A	76	VAL	4

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Mol	Chain	Res	Type	Models (Total)
1	A	69	ALA	4
1	A	39	MET	3
1	A	57	ARG	3
1	A	6	ASP	3
1	A	5	ILE	3
1	A	130	GLY	3
1	A	78	GLN	3
1	A	77	ASP	2
1	A	100	ASN	2
1	A	68	ASN	2
1	A	32	SER	2
1	A	67	GLU	1
1	A	21	ASN	1
1	A	101	CYS	1
1	A	131	THR	1
1	A	41	ILE	1
1	A	20	ASP	1

6.3.2 Protein sidechains [i](#)

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	117/142 (82%)	111±3 (95±3%)	6±3 (5±3%)	31	77
All	All	2340/2840 (82%)	2215 (95%)	125 (5%)	31	77

All 54 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	136	GLU	12
1	A	82	ASN	8
1	A	78	GLN	7
1	A	67	GLU	6
1	A	124	ARG	5
1	A	68	ASN	5
1	A	57	ARG	4
1	A	110	LYS	4

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Mol	Chain	Res	Type	Models (Total)
1	A	119	ASN	3
1	A	134	ASN	3
1	A	47	GLN	3
1	A	72	GLN	3
1	A	37	HIS	3
1	A	84	ARG	3
1	A	22	LYS	3
1	A	115	ASP	2
1	A	21	ASN	2
1	A	103	VAL	2
1	A	120	PHE	2
1	A	30	ASN	2
1	A	77	ASP	2
1	A	62	LEU	2
1	A	58	LEU	2
1	A	18	LEU	2
1	A	42	SER	2
1	A	127	LYS	2
1	A	39	MET	2
1	A	40	THR	2
1	A	109	ASP	2
1	A	137	ASN	1
1	A	29	ARG	1
1	A	46	GLU	1
1	A	8	ARG	1
1	A	128	ILE	1
1	A	19	THR	1
1	A	85	ARG	1
1	A	132	THR	1
1	A	34	MET	1
1	A	28	GLU	1
1	A	139	TYR	1
1	A	114	ASN	1
1	A	116	VAL	1
1	A	118	CYS	1
1	A	66	ASP	1
1	A	31	CYS	1
1	A	100	ASN	1
1	A	23	THR	1
1	A	74	ARG	1
1	A	48	ASN	1
1	A	5	ILE	1

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Mol	Chain	Res	Type	Models (Total)
1	A	59	SER	1
1	A	20	ASP	1
1	A	101	CYS	1
1	A	133	LEU	1

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

6.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 90% for the well-defined parts and 88% for the entire structure.

7.1 Chemical shift list 1

File name: 2ls8_cs.str

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	1889
Number of shifts mapped to atoms	1889
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	13

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	147	-0.02 ± 0.22	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	141	-0.09 ± 0.19	None needed (< 0.5 ppm)
$^{13}\text{C}'$	142	0.07 ± 0.15	None needed (< 0.5 ppm)
^{15}N	131	0.35 ± 0.48	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 90%, i.e. 1512 atoms were assigned a chemical shift out of a possible 1686. 0 out of 15 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	612/638 (96%)	245/254 (96%)	251/260 (97%)	116/124 (94%)
Sidechain	697/835 (83%)	438/494 (89%)	233/291 (80%)	26/50 (52%)

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	Total	¹H	¹³C	¹⁵N
Aromatic	203/213 (95%)	105/111 (95%)	91/91 (100%)	7/11 (64%)
Overall	1512/1686 (90%)	788/859 (92%)	575/642 (90%)	149/185 (81%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 88%, i.e. 1615 atoms were assigned a chemical shift out of a possible 1825. 0 out of 16 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	653/696 (94%)	261/277 (94%)	269/284 (95%)	123/135 (91%)
Sidechain	759/916 (83%)	478/542 (88%)	252/320 (79%)	29/54 (54%)
Aromatic	203/213 (95%)	105/111 (95%)	91/91 (100%)	7/11 (64%)
Overall	1615/1825 (88%)	844/930 (91%)	612/695 (88%)	159/200 (80%)

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	66	ASP	HB3	0.57	4.07 – 1.27	-7.5
1	A	106	TYR	HB2	0.30	4.76 – 1.06	-7.1
1	A	16	PHE	HB3	0.49	4.85 – 1.05	-6.5
1	A	113	TRP	HE3	4.97	9.33 – 5.33	-5.9
1	A	87	PHE	HB3	0.75	4.85 – 1.05	-5.8
1	A	65	ARG	HD2	1.89	4.27 – 1.97	-5.3
1	A	65	ARG	HD3	1.80	4.36 – 1.86	-5.2
1	A	47	GLN	HE21	4.84	9.53 – 4.93	-5.2
1	A	73	TRP	HB3	1.27	4.94 – 1.34	-5.2
1	A	64	LEU	HD21	-0.68	2.14 – -0.66	-5.1
1	A	64	LEU	HD22	-0.68	2.14 – -0.66	-5.1
1	A	64	LEU	HD23	-0.68	2.14 – -0.66	-5.1
1	A	11	GLN	HG3	0.84	3.75 – 0.85	-5.0

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:

