



Full wwPDB NMR Structure Validation Report ⓘ

Apr 27, 2016 – 12:25 AM BST

PDB ID : 2L6Z
Title : haddock model of GATA1NF:Lmo2LIM2-Ldb1LID with FOG
Authors : Wilkinson-White, L.; Gamsjaeger, R.; Dastmalchi, S.; Wienert, B.; Stokes, P.H.; Crossley, M.; Mackay, J.P.; Matthews, J.M.
Deposited on : 2010-12-01

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

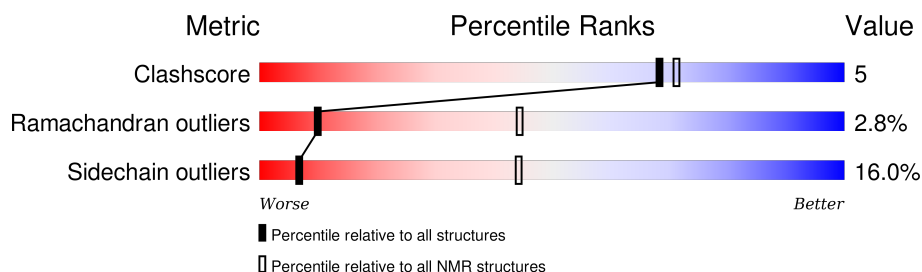
1 Overall quality at a glance

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	39	<div> <div>69%</div> <div>26%</div> <div>5%</div> </div>
2	B	36	<div> <div>72%</div> <div>25%</div> <div>.</div> </div>
3	C	96	<div> <div>70%</div> <div>19%</div> <div>11%</div> </div>

2 Ensemble composition and analysis

This entry contains 10 models. Model 7 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:200-A:238, B:1-B:36, C:91-C:175 (160)	0.45	7

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, 4, 5, 6, 7, 9
2	2, 3, 8
Single-model clusters	10

3 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2589 atoms, of which 1269 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Erythroid transcription factor.

Mol	Chain	Residues	Atoms						Trace
1	A	39	Total	C	H	N	O	S	0
			592	183	286	62	56	5	

- Molecule 2 is a protein called Zinc finger protein ush.

Mol	Chain	Residues	Atoms						Trace
2	B	36	Total	C	H	N	O	S	0
			548	176	272	48	48	4	

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	GLY	-	EXPRESSION TAG	UNP Q9VPQ6
B	2	SER	-	EXPRESSION TAG	UNP Q9VPQ6

- Molecule 3 is a protein called LIM domain only 2, linker, LIM domain-binding protein 1.

Mol	Chain	Residues	Atoms						Trace
3	C	96	Total	C	H	N	O	S	0
			1445	457	711	127	140	10	

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	130	SER	CYS	ENGINEERED MUTATION	UNP Q544Z2

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

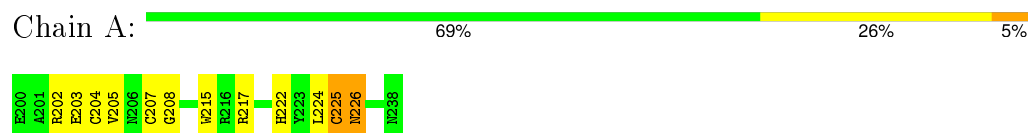
Mol	Chain	Residues	Atoms	
4	B	1	Total	Zn
			1	1
4	A	1	Total	Zn
			1	1
4	C	2	Total	Zn
			2	2

4 Residue-property plots [i](#)

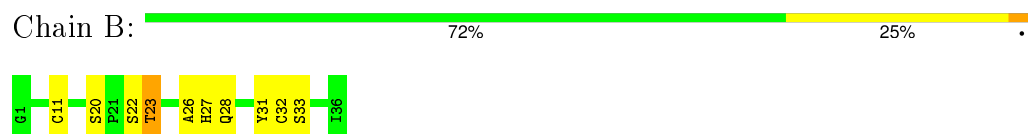
4.1 Average score per residue in the NMR ensemble

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.

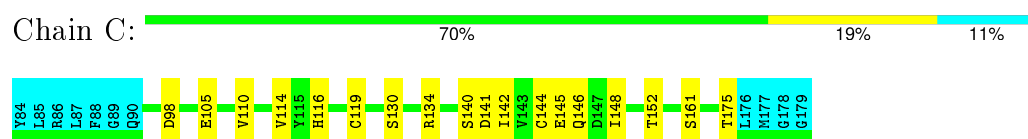
- Molecule 1: Erythroid transcription factor



- Molecule 2: Zinc finger protein ush



- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1

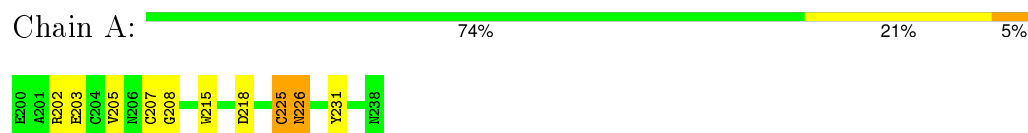


4.2 Scores per residue for each member of the ensemble

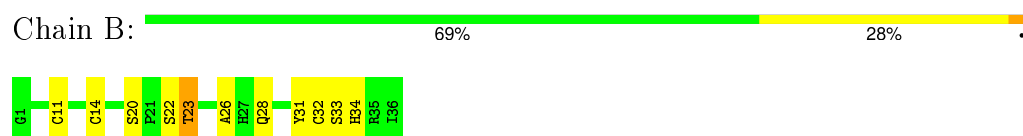
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

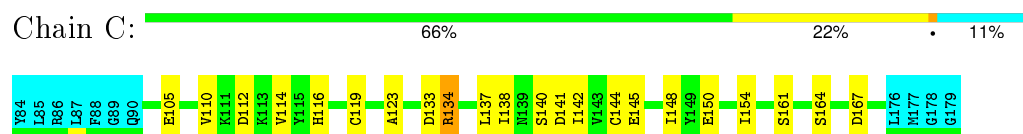
- Molecule 1: Erythroid transcription factor



- Molecule 2: Zinc finger protein ush

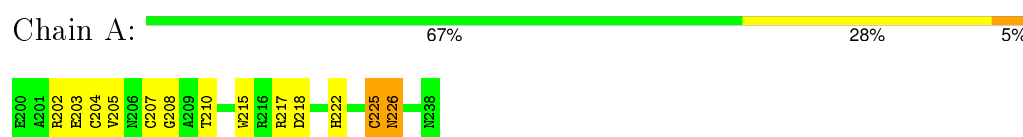


- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1

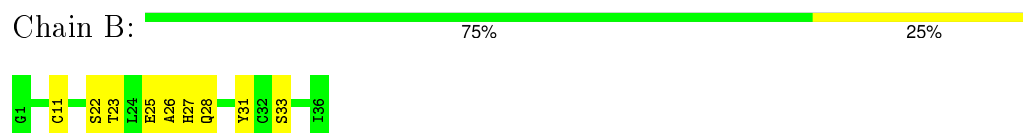


4.2.2 Score per residue for model 2

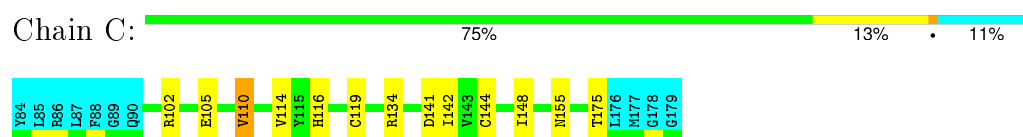
- Molecule 1: Erythroid transcription factor



- Molecule 2: Zinc finger protein ush

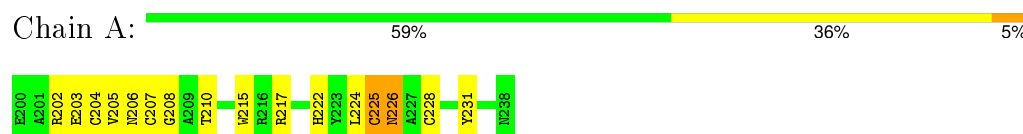


- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1

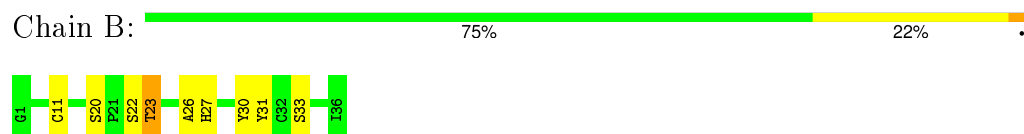


4.2.3 Score per residue for model 3

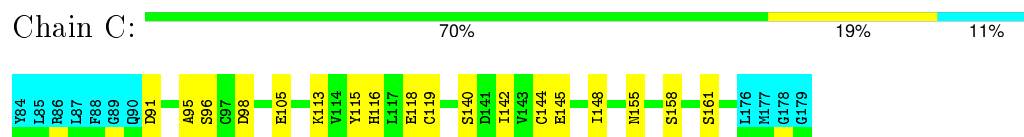
- Molecule 1: Erythroid transcription factor



- Molecule 2: Zinc finger protein ush



- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1

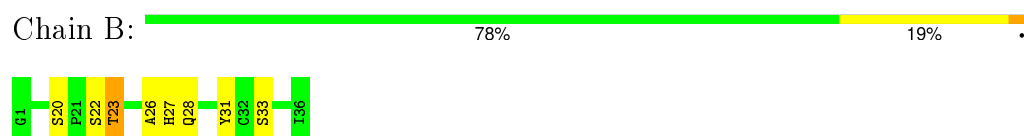


4.2.4 Score per residue for model 4

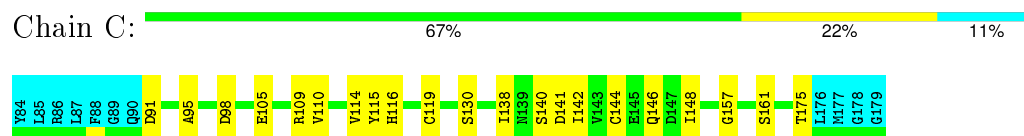
- Molecule 1: Erythroid transcription factor



- Molecule 2: Zinc finger protein ush

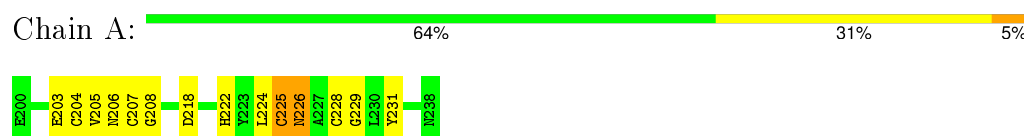


- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1



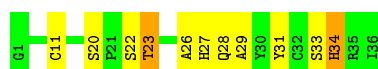
4.2.5 Score per residue for model 5

- Molecule 1: Erythroid transcription factor



- Molecule 2: Zinc finger protein ush





- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1

Chain C: 67% 21% 11%



4.2.6 Score per residue for model 6

- Molecule 1: Erythroid transcription factor

Chain A: 72% 23% 5%



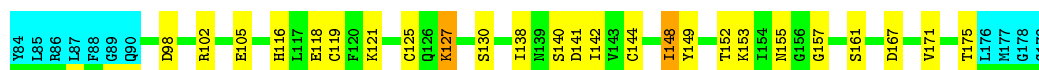
- Molecule 2: Zinc finger protein ush

Chain B: 81% 19%



- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1

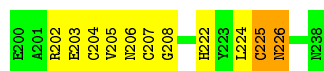
Chain C: 63% 24% 11%



4.2.7 Score per residue for model 7 (medoid)

- Molecule 1: Erythroid transcription factor

Chain A: 72% 23% 5%

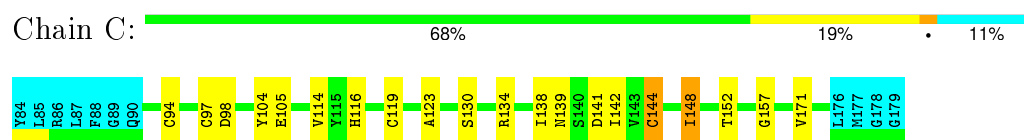


- Molecule 2: Zinc finger protein ush

Chain B: 75% 22%

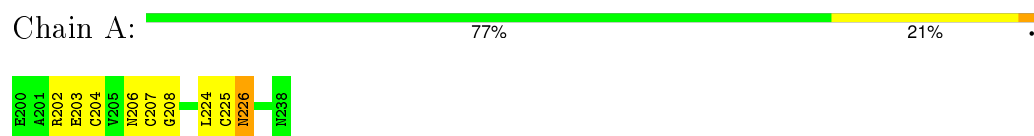


- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1

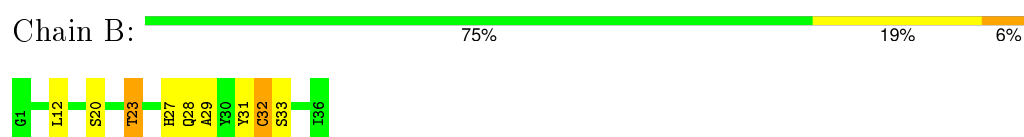


4.2.8 Score per residue for model 8

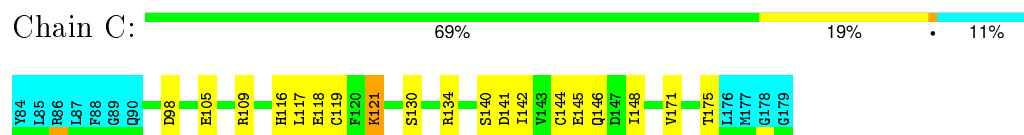
- Molecule 1: Erythroid transcription factor



- Molecule 2: Zinc finger protein ush

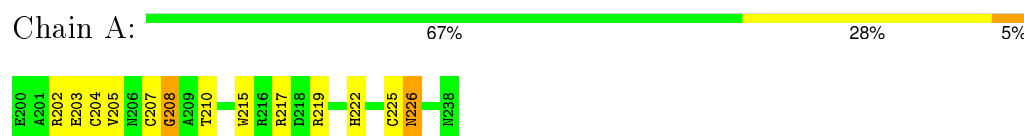


- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1

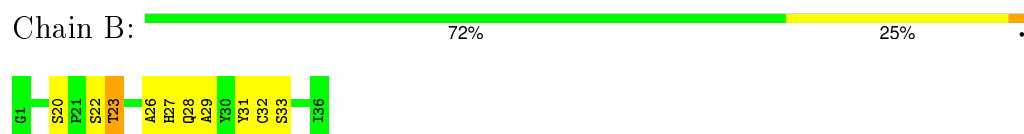


4.2.9 Score per residue for model 9

- Molecule 1: Erythroid transcription factor

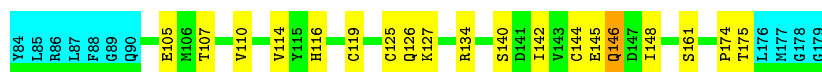


- Molecule 2: Zinc finger protein ush



- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1

Chain C:  69% 19% 11%



4.2.10 Score per residue for model 10

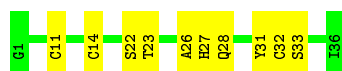
- Molecule 1: Erythroid transcription factor

Chain A:  72% 26%



- Molecule 2: Zinc finger protein ush

Chain B:  72% 28%



- Molecule 3: LIM domain only 2, linker, LIM domain-binding protein 1

Chain C:  72% 16% 11%



5 Refinement protocol and experimental data overview ⓘ

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

Of the 10 calculated structures, 10 were deposited, based on the following criterion: *all calculated structures submitted*.

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

Software name	Classification	Version
CNS	refinement	

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 ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section:
ZN

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	306	286	281	5±1
2	B	276	272	271	3±0
3	C	646	619	618	5±1
All	All	12320	11770	11700	112

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 5.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
3:C:116:HIS:HB2	3:C:119:CYS:SG	0.63	2.34	4	10
3:C:144:CYS:O	3:C:148:ILE:HB	0.62	1.93	4	10
1:A:205:VAL:HB	2:B:26:ALA:HB3	0.60	1.72	5	9
3:C:95:ALA:HB3	3:C:115:TYR:HE1	0.60	1.56	4	2
1:A:207:CYS:SG	1:A:208:GLY:N	0.59	2.75	4	10
2:B:14:CYS:SG	2:B:32:CYS:HA	0.57	2.40	1	2
1:A:225:CYS:SG	1:A:226:ASN:N	0.55	2.79	7	10
1:A:224:LEU:HD22	1:A:229:GLY:HA2	0.54	1.79	5	1
3:C:134:ARG:HG2	3:C:145:GLU:OE1	0.53	2.04	5	1
1:A:204:CYS:HB3	1:A:208:GLY:N	0.53	2.19	9	6

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:20:SER:HB2	2:B:23:THR:OG1	0.52	2.04	7	7
3:C:138:ILE:HG22	3:C:139:ASN:H	0.51	1.65	7	1
3:C:94:CYS:HB3	3:C:97:CYS:O	0.51	2.06	7	1
3:C:95:ALA:HB3	3:C:115:TYR:CE1	0.51	2.40	3	2
3:C:102:ARG:NE	3:C:102:ARG:HA	0.48	2.23	6	2
3:C:150:GLU:O	3:C:154:ILE:HG13	0.47	2.08	1	1
3:C:123:ALA:HB2	3:C:141:ASP:HB2	0.47	1.87	7	2
3:C:144:CYS:SG	3:C:146:GLN:HB3	0.46	2.51	10	4
1:A:204:CYS:SG	1:A:205:VAL:N	0.46	2.88	10	3
1:A:206:ASN:HB3	1:A:228:CYS:SG	0.46	2.50	5	2
3:C:152:THR:HB	3:C:157:GLY:O	0.46	2.11	10	1
2:B:27:HIS:CD2	2:B:27:HIS:C	0.46	2.89	8	3
2:B:27:HIS:C	2:B:27:HIS:CD2	0.46	2.90	10	5
3:C:116:HIS:H	3:C:119:CYS:HB2	0.45	1.71	2	1
1:A:210:THR:HA	1:A:215:TRP:CZ2	0.45	2.47	3	4
3:C:125:CYS:SG	3:C:127:LYS:HG2	0.45	2.52	9	1
3:C:149:TYR:O	3:C:153:LYS:HE2	0.44	2.13	6	1
3:C:121:LYS:HD3	3:C:121:LYS:N	0.44	2.27	8	1
2:B:28:GLN:HA	2:B:32:CYS:SG	0.43	2.54	8	1
3:C:125:CYS:SG	3:C:127:LYS:HB2	0.43	2.54	6	1
3:C:110:VAL:HG12	3:C:111:LYS:H	0.43	1.74	5	1
3:C:130:SER:HB3	3:C:133:ASP:OD1	0.42	2.15	5	1
1:A:202:ARG:HB3	1:A:215:TRP:CE2	0.41	2.50	1	1
3:C:107:THR:O	3:C:174:PRO:HA	0.41	2.16	9	1
3:C:96:SER:CB	3:C:119:CYS:HB3	0.40	2.46	3	1
1:A:205:VAL:HB	2:B:26:ALA:CB	0.40	2.46	6	1
3:C:133:ASP:HA	3:C:134:ARG:NH2	0.40	2.31	1	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	37/39 (95%)	29±1 (78±2%)	7±1 (18±2%)	2±0 (5±1%)	5	27
2	B	34/36 (94%)	25±1 (72±2%)	8±1 (24±4%)	1±1 (4±2%)	6	33

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	85/96 (89%)	71±2 (84±2%)	13±2 (15±2%)	1±1 (1±1%)	20	66
All	All	1560/1710 (91%)	1243 (80%)	274 (18%)	43 (3%)	10	44

All 14 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
2	B	33	SER	10
1	A	226	ASN	10
1	A	225	CYS	6
3	C	157	GLY	3
2	B	29	ALA	3
3	C	138	ILE	2
1	A	208	GLY	2
3	C	156	GLY	1
3	C	110	VAL	1
3	C	104	TYR	1
3	C	91	ASP	1
3	C	144	CYS	1
2	B	34	HIS	1
3	C	112	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	31/31 (100%)	26±1 (85±3%)	5±1 (15±3%)	7	45
2	B	30/30 (100%)	25±1 (83±3%)	5±1 (17±3%)	6	42
3	C	70/78 (90%)	59±2 (84±4%)	11±2 (16±4%)	7	44
All	All	1310/1390 (94%)	1100 (84%)	210 (16%)	7	44

All 52 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	203	GLU	10

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Mol	Chain	Res	Type	Models (Total)
3	C	105	GLU	10
2	B	23	THR	10
2	B	31	TYR	10
3	C	142	ILE	10
2	B	22	SER	8
3	C	114	VAL	7
3	C	140	SER	7
1	A	202	ARG	7
1	A	222	HIS	7
2	B	28	GLN	7
3	C	175	THR	7
2	B	11	CYS	6
3	C	134	ARG	6
1	A	224	LEU	6
3	C	141	ASP	5
1	A	217	ARG	5
3	C	98	ASP	5
3	C	161	SER	5
3	C	110	VAL	5
1	A	231	TYR	4
3	C	130	SER	4
3	C	155	ASN	4
3	C	152	THR	4
3	C	145	GLU	4
1	A	218	ASP	4
3	C	171	VAL	3
2	B	34	HIS	3
3	C	118	GLU	3
2	B	30	TYR	2
3	C	121	LYS	2
3	C	148	ILE	2
2	B	25	GLU	2
3	C	164	SER	2
3	C	117	LEU	2
3	C	146	GLN	2
1	A	206	ASN	2
3	C	109	ARG	2
3	C	167	ASP	2
2	B	32	CYS	2
1	A	200	GLU	1
3	C	138	ILE	1
3	C	113	LYS	1

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Mol	Chain	Res	Type	Models (Total)
3	C	127	LYS	1
3	C	108	MET	1
3	C	126	GLN	1
3	C	91	ASP	1
1	A	219	ARG	1
3	C	137	LEU	1
3	C	158	SER	1
2	B	12	LEU	1
1	A	216	ARG	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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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