



# Full wwPDB NMR Structure Validation Report ⓘ

May 28, 2020 – 10:48 pm BST

PDB ID : 2KXF  
Title : Solution structure of the first two RRM domains of FBP-interacting repressor (FIR)  
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Deposited on : 2010-05-04

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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

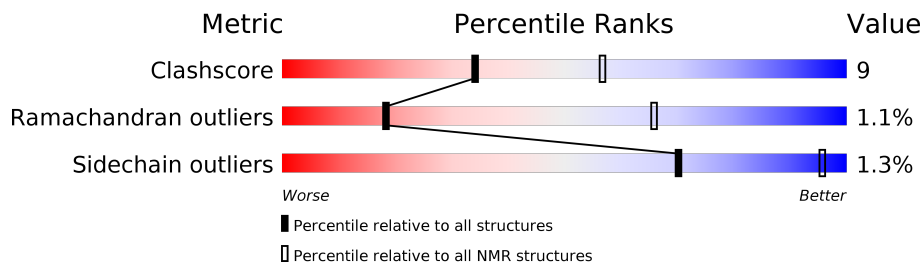
# 1 Overall quality at a glance i

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	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  $\geq 3$ , 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	199	 72% 13% 16%

## 2 Ensemble composition and analysis

This entry contains 20 models. Model 12 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:105-A:144, A:155-A:186, A:193-A:288 (168)	0.52	12

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 3 clusters and 1 single-model cluster was found.

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

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3070 atoms, of which 1532 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Poly(U)-binding-splicing factor PUF60.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	199	3070	973	1532	264	291	10	0

There are 3 discrepancies between the modelled and reference sequences:

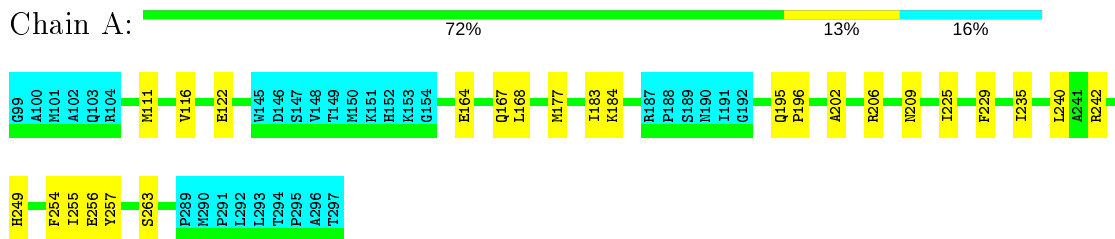
Chain	Residue	Modelled	Actual	Comment	Reference
A	99	GLY	-	expression tag	UNP Q9UHX1
A	100	ALA	-	expression tag	UNP Q9UHX1
A	101	MET	-	expression tag	UNP Q9UHX1

## 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: Poly(U)-binding-splicing factor PUF60

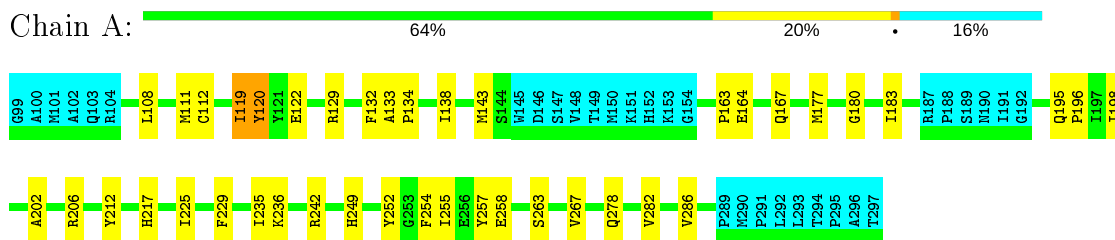


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

#### 4.2.1 Score per residue for model 1

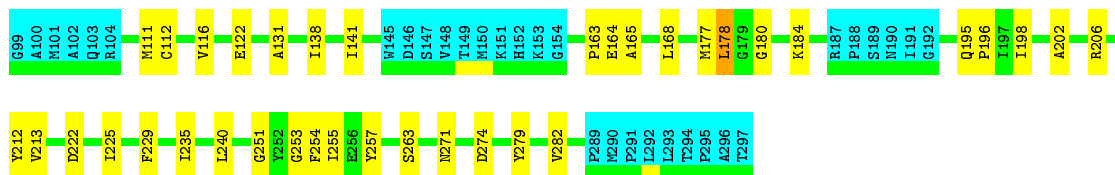
- Molecule 1: Poly(U)-binding-splicing factor PUF60



#### 4.2.2 Score per residue for model 2

- Molecule 1: Poly(U)-binding-splicing factor PUF60

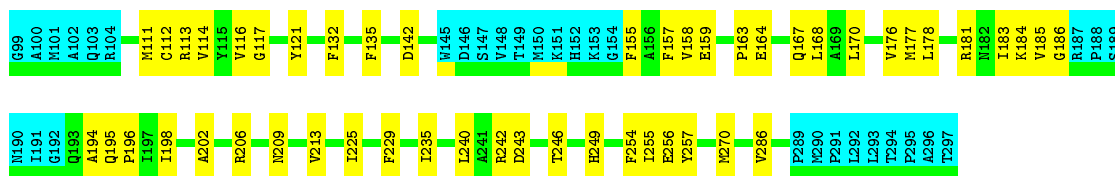




#### 4.2.3 Score per residue for model 3

- Molecule 1: Poly(U)-binding-splicing factor PUF60

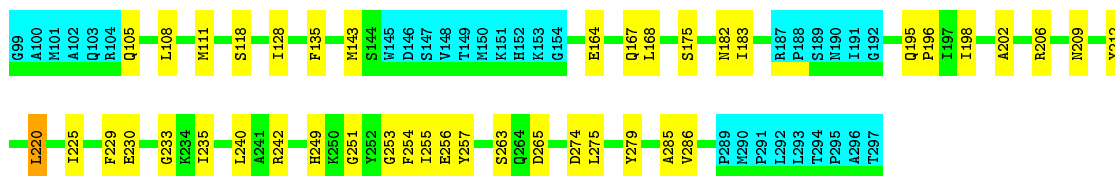
Chain A: 60% 25% 16%



#### 4.2.4 Score per residue for model 4

- Molecule 1: Poly(U)-binding-splicing factor PUF60

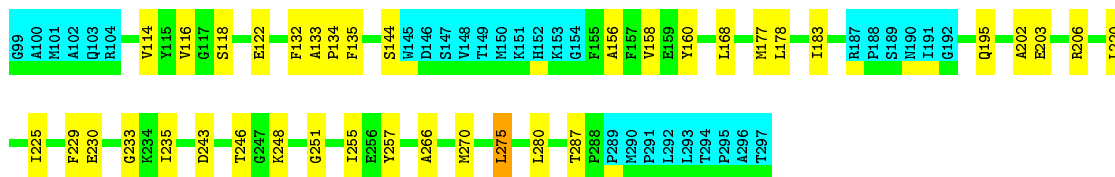
Chain A: 63% 21% 16%



#### 4.2.5 Score per residue for model 5

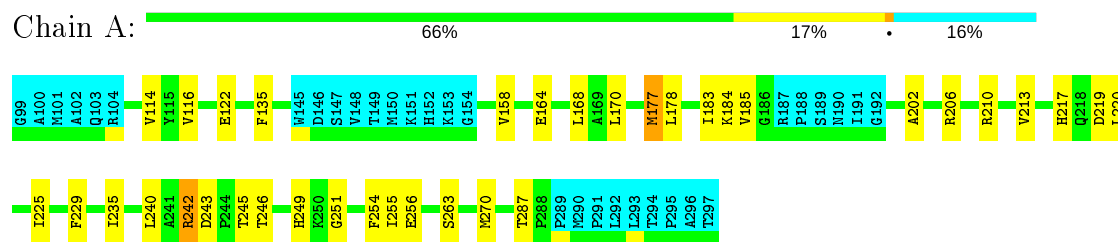
- Molecule 1: Poly(U)-binding-splicing factor PUF60

Chain A: 66% 18% 16%



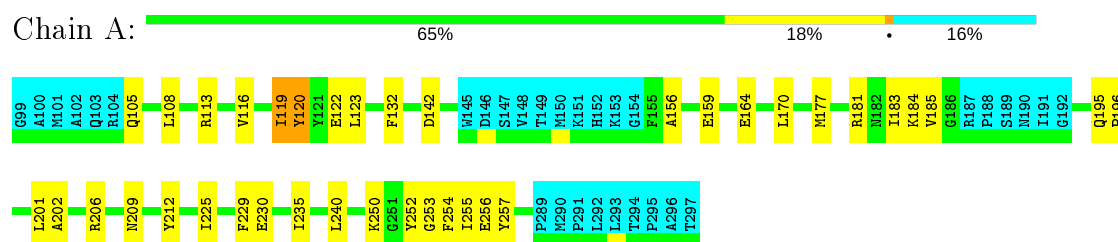
#### 4.2.6 Score per residue for model 6

- Molecule 1: Poly(U)-binding-splicing factor PUF60



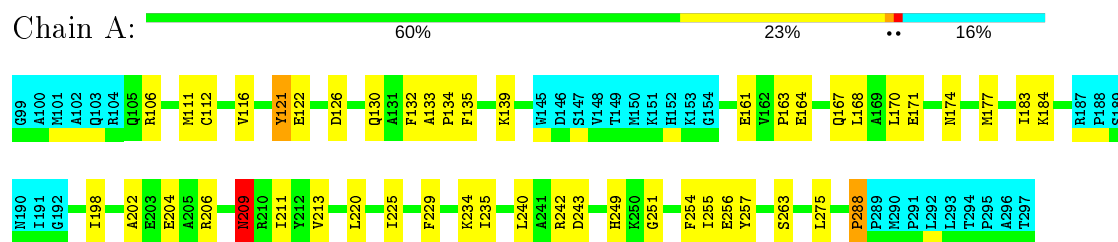
#### 4.2.7 Score per residue for model 7

- Molecule 1: Poly(U)-binding-splicing factor PUF60



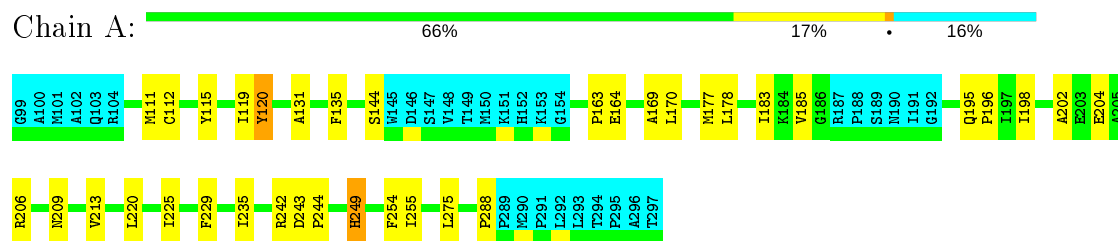
#### 4.2.8 Score per residue for model 8

- Molecule 1: Poly(U)-binding-splicing factor PUF60



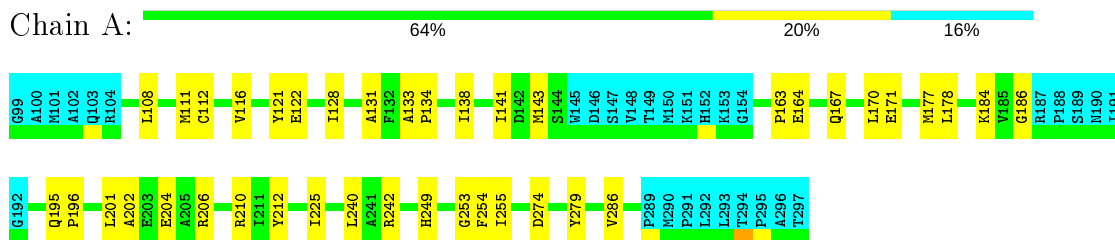
#### 4.2.9 Score per residue for model 9

- Molecule 1: Poly(U)-binding-splicing factor PUF60



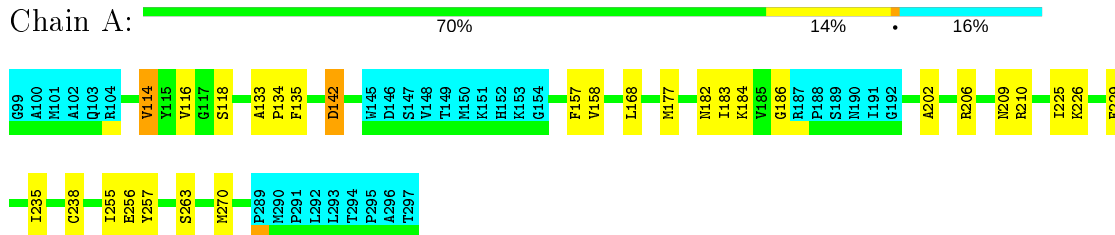
#### 4.2.10 Score per residue for model 10

- Molecule 1: Poly(U)-binding-splicing factor PUF60



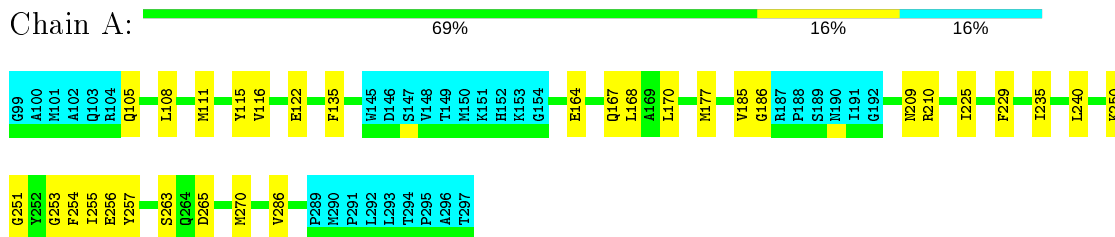
#### 4.2.11 Score per residue for model 11

- Molecule 1: Poly(U)-binding-splicing factor PUF60



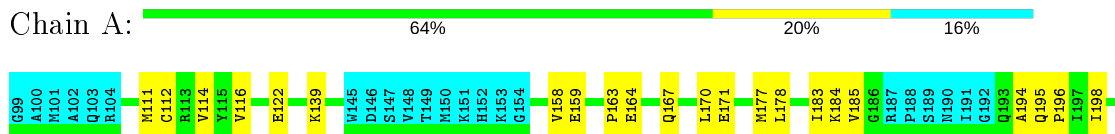
#### 4.2.12 Score per residue for model 12 (medoid)

- Molecule 1: Poly(U)-binding-splicing factor PUF60



#### 4.2.13 Score per residue for model 13

- Molecule 1: Poly(U)-binding-splicing factor PUF60







#### 4.2.14 Score per residue for model 14

- Molecule 1: Poly(U)-binding-splicing factor PUF60

Chain A: 65% 19% 16%



#### 4.2.15 Score per residue for model 15

- Molecule 1: Poly(U)-binding-splicing factor PUF60

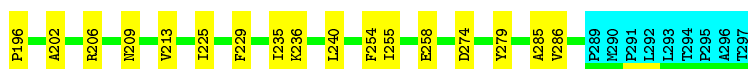
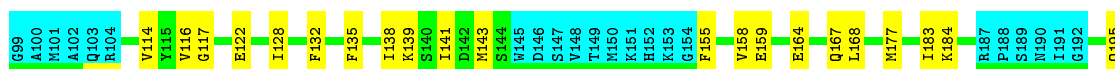
Chain A: 64% 20% 16%



#### 4.2.16 Score per residue for model 16

- Molecule 1: Poly(U)-binding-splicing factor PUF60

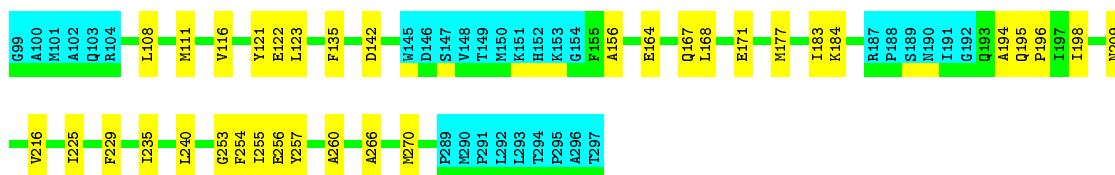
Chain A: 65% 19% 16%



#### 4.2.17 Score per residue for model 17

- Molecule 1: Poly(U)-binding-splicing factor PUF60

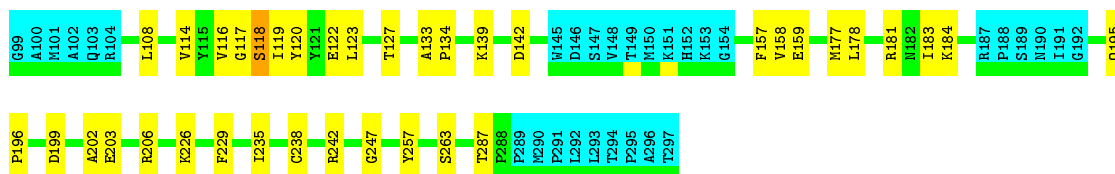
Chain A: 67% 17% 16%



#### 4.2.18 Score per residue for model 18

- Molecule 1: Poly(U)-binding-splicing factor PUF60

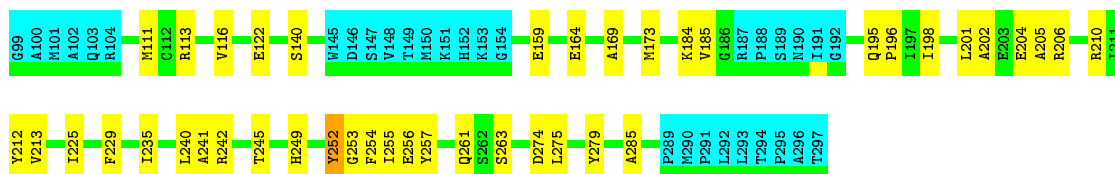
Chain A: 66% 18% 16%



#### 4.2.19 Score per residue for model 19

- Molecule 1: Poly(U)-binding-splicing factor PUF60

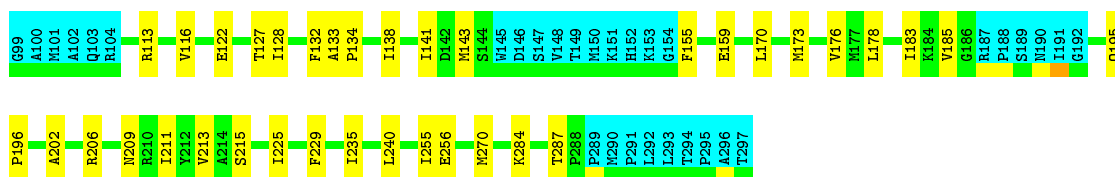
Chain A: 63% 21% 16%



#### 4.2.20 Score per residue for model 20

- Molecule 1: Poly(U)-binding-splicing factor PUF60

Chain A: 66% 18% 16%



## 5 Refinement protocol and experimental data overview

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

Of the 200 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
ARIA	refinement	1.2
ARIA	structure solution	1.2

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 i

### 6.1 Standard geometry i

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 i

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	1305	1290	1288	23±3
All	All	26100	25800	25760	460

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 9.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:177:MET:HB2	1:A:183:ILE:HD13	0.81	1.51	15	13
1:A:210:ARG:HB3	1:A:256:GLU:HG2	0.79	1.53	14	4
1:A:112:CYS:HA	1:A:163:PRO:HA	0.76	1.55	10	8
1:A:113:ARG:HG2	1:A:159:GLU:HG3	0.74	1.59	19	2
1:A:177:MET:HE2	1:A:178:LEU:HG	0.73	1.60	13	4
1:A:113:ARG:HG2	1:A:159:GLU:HG2	0.71	1.59	7	1
1:A:114:VAL:HG12	1:A:158:VAL:HB	0.71	1.62	3	7
1:A:225:ILE:HG22	1:A:255:ILE:HD11	0.70	1.64	1	19
1:A:164:GLU:HG3	1:A:254:PHE:HE1	0.69	1.47	2	1
1:A:164:GLU:HG3	1:A:254:PHE:HE2	0.69	1.47	13	1
1:A:140:SER:HB3	1:A:159:GLU:HB2	0.68	1.65	19	1
1:A:118:SER:HB3	1:A:183:ILE:HG22	0.64	1.68	18	1
1:A:116:VAL:HA	1:A:184:LYS:O	0.63	1.93	3	12
1:A:118:SER:HB2	1:A:183:ILE:HA	0.63	1.69	11	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:138:ILE:HD13	1:A:141:ILE:HD11	0.61	1.72	10	4
1:A:229:PHE:HB2	1:A:235:ILE:HD11	0.61	1.72	17	18
1:A:117:GLY:HA3	1:A:155:PHE:HD1	0.61	1.56	3	2
1:A:202:ALA:O	1:A:206:ARG:HG3	0.58	1.99	5	18
1:A:242:ARG:HG2	1:A:249:HIS:HB3	0.57	1.76	10	2
1:A:225:ILE:HG13	1:A:240:LEU:HD21	0.57	1.76	8	6
1:A:142:ASP:HB2	1:A:157:PHE:HB2	0.57	1.77	3	4
1:A:226:LYS:HD2	1:A:238:CYS:SG	0.56	2.40	11	1
1:A:169:ALA:HB1	1:A:173:MET:SD	0.56	2.41	19	1
1:A:133:ALA:N	1:A:134:PRO:HD2	0.55	2.15	1	8
1:A:177:MET:HA	1:A:177:MET:HE3	0.55	1.79	3	2
1:A:184:LYS:HE2	1:A:184:LYS:HA	0.55	1.77	19	1
1:A:226:LYS:HB2	1:A:238:CYS:HB3	0.55	1.79	18	1
1:A:242:ARG:HG3	1:A:249:HIS:HB3	0.55	1.78	6	1
1:A:195:GLN:HB2	1:A:196:PRO:HD3	0.55	1.78	1	4
1:A:210:ARG:HB3	1:A:256:GLU:CG	0.55	2.31	6	1
1:A:220:LEU:HD21	1:A:275:LEU:HG	0.55	1.77	4	3
1:A:108:LEU:HD21	1:A:287:THR:HG21	0.55	1.79	18	1
1:A:242:ARG:HA	1:A:249:HIS:HA	0.54	1.77	10	6
1:A:274:ASP:HB3	1:A:279:TYR:CE1	0.54	2.38	4	3
1:A:167:GLN:HG3	1:A:286:VAL:HG11	0.54	1.78	15	6
1:A:173:MET:SD	1:A:183:ILE:HD11	0.54	2.42	20	1
1:A:164:GLU:HG3	1:A:254:PHE:CE1	0.54	2.38	9	3
1:A:217:HIS:HB3	1:A:220:LEU:HD23	0.54	1.79	6	1
1:A:167:GLN:HG2	1:A:286:VAL:HG13	0.54	1.80	16	1
1:A:135:PHE:HD1	1:A:168:LEU:HG	0.54	1.63	17	8
1:A:119:ILE:H	1:A:183:ILE:HG22	0.54	1.63	18	1
1:A:195:GLN:HB3	1:A:196:PRO:HD3	0.53	1.80	19	3
1:A:135:PHE:HD2	1:A:168:LEU:HG	0.53	1.63	8	1
1:A:118:SER:HB3	1:A:183:ILE:HA	0.53	1.79	15	2
1:A:257:TYR:HB2	1:A:263:SER:OG	0.53	2.04	19	2
1:A:217:HIS:HA	1:A:278:GLN:OE1	0.53	2.04	1	1
1:A:195:GLN:N	1:A:196:PRO:HD2	0.52	2.20	20	8
1:A:170:LEU:HD13	1:A:185:VAL:O	0.52	2.04	13	3
1:A:204:GLU:HB3	1:A:288:PRO:HG3	0.52	1.81	9	2
1:A:111:MET:SD	1:A:198:ILE:HG12	0.52	2.44	1	7
1:A:142:ASP:O	1:A:156:ALA:HA	0.52	2.04	7	2
1:A:123:LEU:HD13	1:A:177:MET:HG3	0.52	1.81	17	1
1:A:117:GLY:HA3	1:A:155:PHE:CD1	0.52	2.38	16	2
1:A:218:GLN:HA	1:A:249:HIS:CE1	0.52	2.39	15	1
1:A:108:LEU:HD21	1:A:201:LEU:HD13	0.52	1.81	7	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:164:GLU:HG2	1:A:254:PHE:HE2	0.52	1.65	4	5
1:A:209:ASN:O	1:A:256:GLU:HA	0.51	2.05	17	10
1:A:210:ARG:CB	1:A:256:GLU:HG2	0.51	2.32	14	2
1:A:257:TYR:CD1	1:A:263:SER:HA	0.51	2.41	11	2
1:A:230:GLU:HA	1:A:233:GLY:O	0.51	2.04	5	2
1:A:114:VAL:CG1	1:A:158:VAL:HB	0.51	2.33	3	1
1:A:138:ILE:HG21	1:A:141:ILE:HG13	0.51	1.81	10	1
1:A:111:MET:SD	1:A:198:ILE:HG23	0.51	2.46	4	1
1:A:173:MET:HG2	1:A:176:VAL:O	0.50	2.06	15	1
1:A:117:GLY:O	1:A:118:SER:HB2	0.50	2.06	18	1
1:A:243:ASP:HB2	1:A:246:THR:OG1	0.50	2.06	6	3
1:A:203:GLU:OE1	1:A:206:ARG:HD3	0.50	2.07	5	1
1:A:211:ILE:HG22	1:A:284:LYS:HD2	0.50	1.83	20	1
1:A:229:PHE:CB	1:A:235:ILE:HD11	0.50	2.36	17	15
1:A:164:GLU:HG3	1:A:254:PHE:CE2	0.49	2.36	13	2
1:A:170:LEU:O	1:A:174:ASN:HB2	0.49	2.07	15	1
1:A:212:TYR:HD1	1:A:254:PHE:CE1	0.49	2.24	7	3
1:A:167:GLN:HB2	1:A:286:VAL:HG11	0.49	1.83	10	1
1:A:170:LEU:HD12	1:A:186:GLY:HA3	0.49	1.83	3	2
1:A:213:VAL:HG11	1:A:229:PHE:CZ	0.49	2.43	19	7
1:A:238:CYS:SG	1:A:255:ILE:HG12	0.49	2.47	15	1
1:A:258:GLU:HB2	1:A:259:LYS:HD2	0.49	1.84	15	1
1:A:242:ARG:HB3	1:A:249:HIS:CB	0.48	2.38	4	1
1:A:177:MET:HE3	1:A:177:MET:HA	0.48	1.85	14	1
1:A:229:PHE:CE2	1:A:270:MET:HG3	0.48	2.43	6	5
1:A:105:GLN:HG3	1:A:108:LEU:HD12	0.48	1.83	12	1
1:A:113:ARG:HG2	1:A:159:GLU:HB2	0.48	1.83	14	3
1:A:257:TYR:HB2	1:A:263:SER:HB3	0.48	1.86	18	6
1:A:164:GLU:HB3	1:A:212:TYR:CD2	0.48	2.43	4	2
1:A:199:ASP:O	1:A:203:GLU:HG3	0.48	2.09	14	2
1:A:139:LYS:HD2	1:A:159:GLU:OE2	0.48	2.09	16	1
1:A:220:LEU:HD21	1:A:275:LEU:HD13	0.48	1.84	5	1
1:A:175:SER:O	1:A:182:ASN:HA	0.47	2.09	4	1
1:A:236:LYS:HB2	1:A:258:GLU:HG2	0.47	1.86	16	2
1:A:164:GLU:HG3	1:A:254:PHE:CZ	0.47	2.44	6	1
1:A:212:TYR:HD1	1:A:254:PHE:CE2	0.47	2.28	10	1
1:A:114:VAL:HG22	1:A:186:GLY:H	0.47	1.70	11	1
1:A:217:HIS:HB3	1:A:220:LEU:CD2	0.47	2.39	6	1
1:A:119:ILE:O	1:A:120:TYR:HB2	0.46	2.10	7	2
1:A:229:PHE:HB3	1:A:257:TYR:OH	0.46	2.10	19	11
1:A:226:LYS:HB2	1:A:238:CYS:SG	0.46	2.50	11	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:241:ALA:HB3	1:A:252:TYR:CD1	0.46	2.45	19	1
1:A:165:ALA:HA	1:A:168:LEU:HB3	0.46	1.86	2	1
1:A:116:VAL:HG11	1:A:132:PHE:CZ	0.46	2.45	20	4
1:A:120:TYR:HA	1:A:181:ARG:HD3	0.46	1.85	18	1
1:A:164:GLU:HG2	1:A:254:PHE:HE1	0.46	1.70	17	3
1:A:209:ASN:OD1	1:A:260:ALA:HA	0.46	2.11	17	1
1:A:203:GLU:HA	1:A:206:ARG:HD3	0.46	1.87	18	1
1:A:132:PHE:HB2	1:A:138:ILE:HD11	0.46	1.87	14	2
1:A:240:LEU:HD23	1:A:253:GLY:HA3	0.46	1.87	4	2
1:A:116:VAL:HG12	1:A:185:VAL:HG12	0.46	1.87	3	3
1:A:119:ILE:O	1:A:181:ARG:HG2	0.46	2.10	7	1
1:A:240:LEU:CD2	1:A:253:GLY:HA3	0.46	2.41	17	2
1:A:271:ASN:HA	1:A:282:VAL:HG12	0.46	1.87	2	1
1:A:240:LEU:HA	1:A:253:GLY:CA	0.46	2.41	14	4
1:A:225:ILE:CG2	1:A:255:ILE:HD11	0.46	2.41	17	9
1:A:111:MET:CG	1:A:198:ILE:HD12	0.46	2.41	9	1
1:A:275:LEU:HD12	1:A:280:LEU:HD11	0.46	1.87	5	1
1:A:135:PHE:CD2	1:A:168:LEU:HG	0.45	2.45	8	1
1:A:105:GLN:HB2	1:A:108:LEU:HD12	0.45	1.88	4	1
1:A:230:GLU:OE1	1:A:235:ILE:HB	0.45	2.11	7	1
1:A:240:LEU:HA	1:A:253:GLY:HA3	0.45	1.87	7	3
1:A:164:GLU:O	1:A:167:GLN:HG3	0.45	2.12	8	2
1:A:167:GLN:CB	1:A:286:VAL:HG11	0.45	2.42	10	1
1:A:108:LEU:HD23	1:A:111:MET:SD	0.45	2.52	1	3
1:A:267:VAL:HA	1:A:282:VAL:CG1	0.45	2.42	1	1
1:A:128:ILE:HG13	1:A:143:MET:CE	0.45	2.42	16	3
1:A:205:ALA:HB2	1:A:285:ALA:O	0.45	2.12	19	1
1:A:164:GLU:HG2	1:A:254:PHE:CE1	0.44	2.46	3	2
1:A:137:PRO:HG3	1:A:252:TYR:CE1	0.44	2.47	15	1
1:A:115:TYR:O	1:A:185:VAL:HA	0.44	2.12	12	2
1:A:221:SER:O	1:A:225:ILE:HG12	0.44	2.12	15	1
1:A:131:ALA:HB1	1:A:177:MET:HE1	0.44	1.88	2	2
1:A:213:VAL:HG12	1:A:282:VAL:HG23	0.44	1.90	13	2
1:A:167:GLN:O	1:A:171:GLU:HG2	0.44	2.13	8	3
1:A:164:GLU:HG2	1:A:254:PHE:CE2	0.44	2.48	12	1
1:A:242:ARG:CB	1:A:249:HIS:HA	0.44	2.42	19	2
1:A:194:ALA:O	1:A:198:ILE:HG13	0.44	2.12	3	4
1:A:139:LYS:HB3	1:A:159:GLU:HG2	0.44	1.88	18	1
1:A:126:ASP:O	1:A:130:GLN:HG3	0.44	2.13	8	1
1:A:262:SER:HA	1:A:265:ASP:OD2	0.44	2.13	14	1
1:A:274:ASP:HB2	1:A:279:TYR:CE1	0.44	2.47	2	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:119:ILE:O	1:A:120:TYR:HB3	0.43	2.13	9	1
1:A:201:LEU:HA	1:A:204:GLU:HB2	0.43	1.90	10	1
1:A:132:PHE:HB3	1:A:160:TYR:OH	0.43	2.13	5	1
1:A:128:ILE:HG23	1:A:132:PHE:CE1	0.43	2.48	20	1
1:A:128:ILE:HG13	1:A:143:MET:HE1	0.43	1.90	4	1
1:A:201:LEU:HA	1:A:204:GLU:HG3	0.43	1.90	19	1
1:A:170:LEU:HA	1:A:185:VAL:CG2	0.43	2.43	12	1
1:A:116:VAL:HG11	1:A:132:PHE:HZ	0.43	1.73	20	1
1:A:170:LEU:HG	1:A:185:VAL:HG23	0.43	1.90	3	2
1:A:226:LYS:HB2	1:A:238:CYS:CB	0.43	2.44	18	1
1:A:211:ILE:HG23	1:A:263:SER:HB2	0.43	1.91	8	2
1:A:242:ARG:HD2	1:A:247:GLY:O	0.43	2.13	18	1
1:A:225:ILE:O	1:A:229:PHE:HD1	0.43	1.97	20	1
1:A:170:LEU:O	1:A:174:ASN:HB3	0.42	2.14	8	1
1:A:164:GLU:HG3	1:A:285:ALA:HA	0.42	1.91	4	2
1:A:167:GLN:HG3	1:A:286:VAL:HG21	0.42	1.91	14	1
1:A:175:SER:HA	1:A:183:ILE:O	0.42	2.14	15	1
1:A:177:MET:SD	1:A:178:LEU:HG	0.42	2.54	18	1
1:A:131:ALA:HB1	1:A:177:MET:SD	0.42	2.55	10	1
1:A:216:VAL:HG13	1:A:225:ILE:HD11	0.42	1.90	17	1
1:A:177:MET:HB3	1:A:183:ILE:HG21	0.42	1.91	5	1
1:A:233:GLY:HA3	1:A:257:TYR:HE2	0.42	1.75	15	1
1:A:164:GLU:HG3	1:A:254:PHE:HZ	0.41	1.75	6	1
1:A:176:VAL:HA	1:A:181:ARG:O	0.41	2.16	3	1
1:A:135:PHE:CD1	1:A:169:ALA:HA	0.41	2.50	9	1
1:A:116:VAL:HG23	1:A:156:ALA:HB3	0.41	1.92	5	1
1:A:266:ALA:O	1:A:270:MET:HG2	0.41	2.15	17	2
1:A:170:LEU:HD13	1:A:186:GLY:HA3	0.41	1.91	10	1
1:A:262:SER:O	1:A:265:ASP:HB2	0.41	2.15	14	1
1:A:201:LEU:O	1:A:204:GLU:HB2	0.41	2.15	19	1
1:A:116:VAL:O	1:A:155:PHE:HA	0.41	2.16	20	1
1:A:212:TYR:HB2	1:A:254:PHE:CE2	0.41	2.51	19	1
1:A:170:LEU:HG	1:A:185:VAL:HG13	0.41	1.93	20	1
1:A:250:LYS:HB2	1:A:252:TYR:CE1	0.41	2.51	7	1
1:A:170:LEU:HA	1:A:185:VAL:HG21	0.41	1.92	9	1
1:A:168:LEU:HD11	1:A:212:TYR:CD2	0.41	2.50	15	1
1:A:177:MET:CB	1:A:183:ILE:HD13	0.41	2.46	16	2
1:A:241:ALA:HB3	1:A:252:TYR:CE1	0.41	2.51	19	1
1:A:263:SER:O	1:A:267:VAL:HG23	0.40	2.16	1	1
1:A:243:ASP:HA	1:A:244:PRO:HD3	0.40	1.82	9	1
1:A:177:MET:O	1:A:178:LEU:HB2	0.40	2.17	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:139:LYS:HE3	1:A:159:GLU:OE1	0.40	2.16	13	1
1:A:129:ARG:HA	1:A:138:ILE:HD12	0.40	1.93	1	1
1:A:138:ILE:CG2	1:A:141:ILE:HG13	0.40	2.47	10	1
1:A:123:LEU:HD22	1:A:127:THR:HG21	0.40	1.93	18	1
1:A:139:LYS:HD2	1:A:161:GLU:HG2	0.40	1.93	8	1
1:A:128:ILE:HG23	1:A:132:PHE:CD1	0.40	2.52	16	1

## 6.3 Torsion angles [i](#)

### 6.3.1 Protein backbone [i](#)

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	168/199 (84%)	158±2 (94±1%)	8±2 (5±1%)	2±1 (1±1%)	18	66
All	All	3360/3980 (84%)	3156 (94%)	166 (5%)	38 (1%)	18	66

All 12 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	122	GLU	14
1	A	251	GLY	7
1	A	120	TYR	3
1	A	178	LEU	3
1	A	180	GLY	2
1	A	209	ASN	2
1	A	119	ILE	2
1	A	250	LYS	1
1	A	245	THR	1
1	A	118	SER	1
1	A	288	PRO	1
1	A	123	LEU	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	137/162 (85%)	135±1 (99±1%)	2±1 (1±1%)	70	96
All	All	2740/3240 (85%)	2704 (99%)	36 (1%)	70	96

All 30 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	177	MET	2
1	A	265	ASP	2
1	A	275	LEU	2
1	A	209	ASN	2
1	A	222	ASP	2
1	A	252	TYR	2
1	A	178	LEU	1
1	A	105	GLN	1
1	A	176	VAL	1
1	A	111	MET	1
1	A	142	ASP	1
1	A	122	GLU	1
1	A	114	VAL	1
1	A	219	ASP	1
1	A	263	SER	1
1	A	234	LYS	1
1	A	121	TYR	1
1	A	215	SER	1
1	A	143	MET	1
1	A	248	LYS	1
1	A	245	THR	1
1	A	261	GLN	1
1	A	171	GLU	1
1	A	243	ASP	1
1	A	127	THR	1
1	A	182	ASN	1
1	A	249	HIS	1
1	A	242	ARG	1
1	A	164	GLU	1

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Mol	Chain	Res	Type	Models (Total)
1	A	220	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

No chemical shift data were provided