



# Full wwPDB NMR Structure Validation Report ⓘ

Feb 12, 2017 – 09:50 pm GMT

PDB ID : 2F3J  
Title : The solution structure of the REF2-I mRNA export factor (residues 1-155).  
Authors : Golovanov, A.P.; Hautbergue, G.M.; Wilson, S.A.; Lian, L.Y.  
Deposited on : 2005-11-21

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  
<http://wwpdb.org/validation/2016/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 : 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

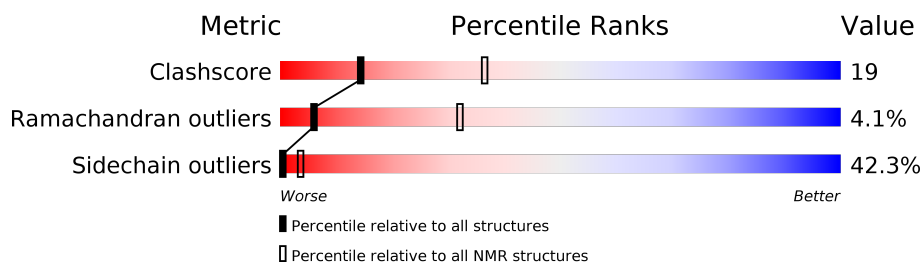
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment 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	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  $\geq 3$ , 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	177	

## 2 Ensemble composition and analysis

This entry contains 14 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:8-A:18 (11)	0.22	3
2	A:44-A:51 (8)	0.38	4
3	A:74-A:151 (78)	0.88	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 3 clusters and 1 single-model cluster was found.

Cluster number	Models
1	5, 6, 7, 8, 11, 12, 13, 14
2	1, 4, 10
3	2, 9
Single-model clusters	3

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2444 atoms, of which 1219 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called RNA and export factor binding protein 2.

Mol	Chain	Residues	Atoms						Trace
1	A	158	Total	C	H	N	O	S	0
			2444	754	1219	235	230	6	

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	CLONING ARTIFACT	UNP Q9JJW6
A	-12	ALA	-	CLONING ARTIFACT	UNP Q9JJW6
A	-11	SER	-	CLONING ARTIFACT	UNP Q9JJW6
A	-10	MET	-	CLONING ARTIFACT	UNP Q9JJW6
A	-9	THR	-	CLONING ARTIFACT	UNP Q9JJW6
A	-8	GLY	-	CLONING ARTIFACT	UNP Q9JJW6
A	-7	GLY	-	CLONING ARTIFACT	UNP Q9JJW6
A	-6	GLN	-	CLONING ARTIFACT	UNP Q9JJW6
A	-5	GLN	-	CLONING ARTIFACT	UNP Q9JJW6
A	-4	MET	-	CLONING ARTIFACT	UNP Q9JJW6
A	-3	GLY	-	CLONING ARTIFACT	UNP Q9JJW6
A	-2	ARG	-	CLONING ARTIFACT	UNP Q9JJW6
A	-1	ASP	-	CLONING ARTIFACT	UNP Q9JJW6
A	0	PRO	-	CLONING ARTIFACT	UNP Q9JJW6
A	156	LEU	-	CLONING ARTIFACT	UNP Q9JJW6
A	157	GLU	-	CLONING ARTIFACT	UNP Q9JJW6
A	158	HIS	-	EXPRESSION TAG	UNP Q9JJW6
A	159	HIS	-	EXPRESSION TAG	UNP Q9JJW6
A	160	HIS	-	EXPRESSION TAG	UNP Q9JJW6
A	161	HIS	-	EXPRESSION TAG	UNP Q9JJW6
A	162	HIS	-	EXPRESSION TAG	UNP Q9JJW6
A	163	HIS	-	EXPRESSION TAG	UNP Q9JJW6

## 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: RNA and export factor binding protein 2

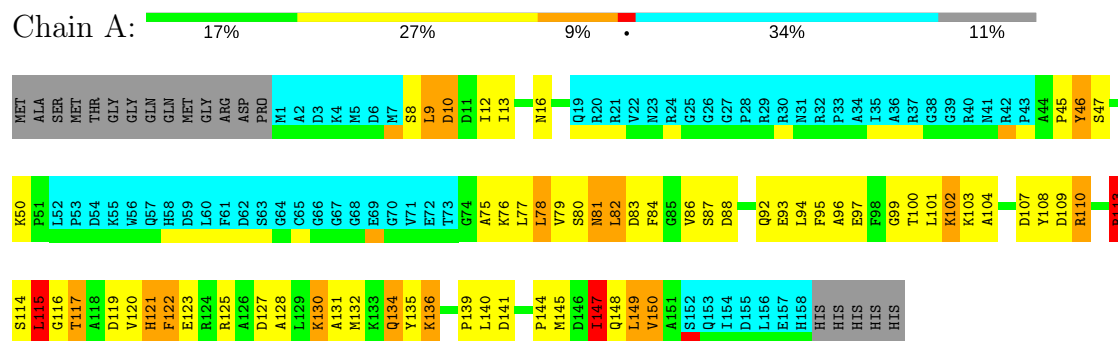


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

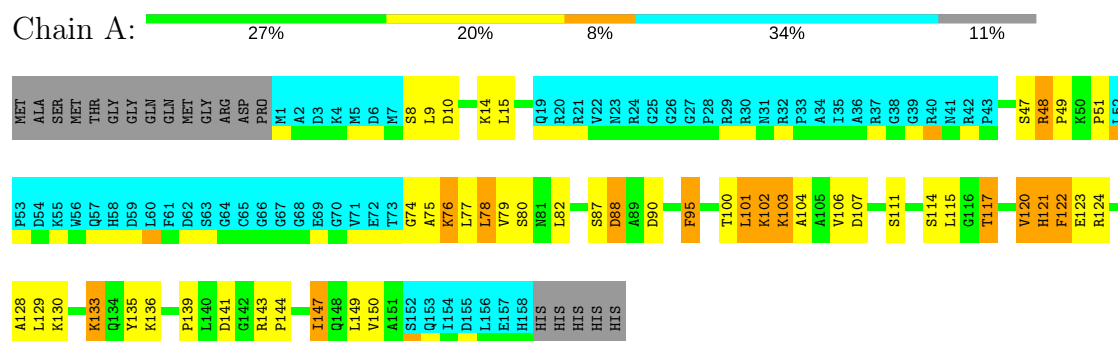
#### 4.2.1 Score per residue for model 1

- Molecule 1: RNA and export factor binding protein 2



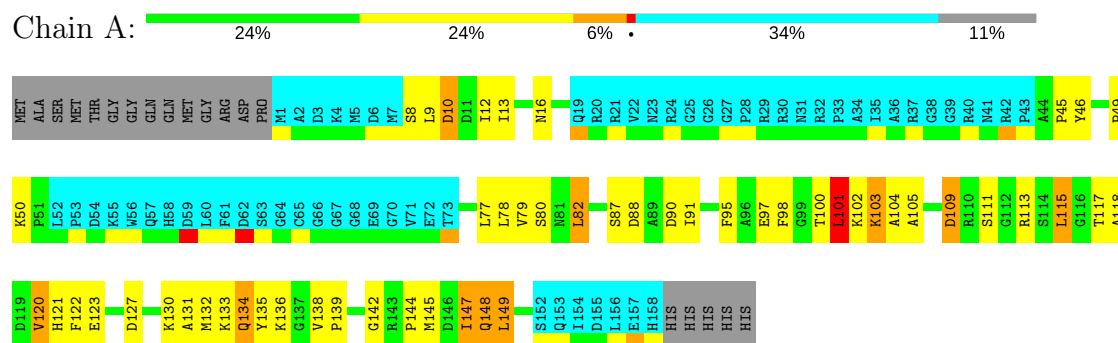
### 4.2.2 Score per residue for model 2

- Molecule 1: RNA and export factor binding protein 2



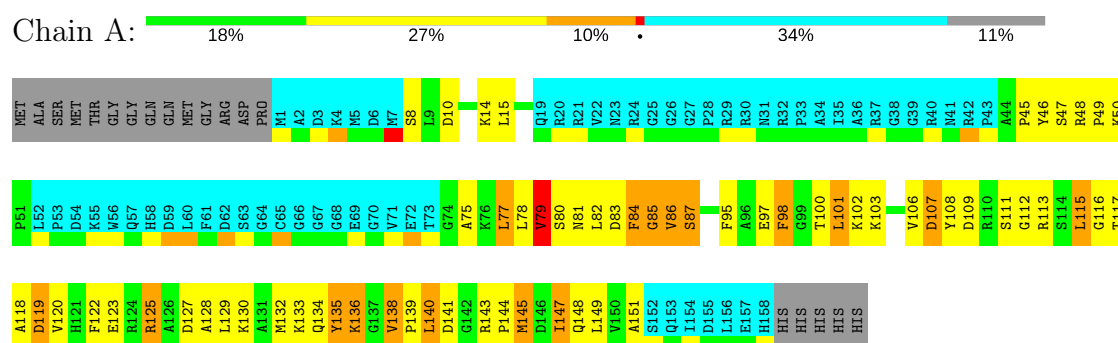
### 4.2.3 Score per residue for model 3

- Molecule 1: RNA and export factor binding protein 2



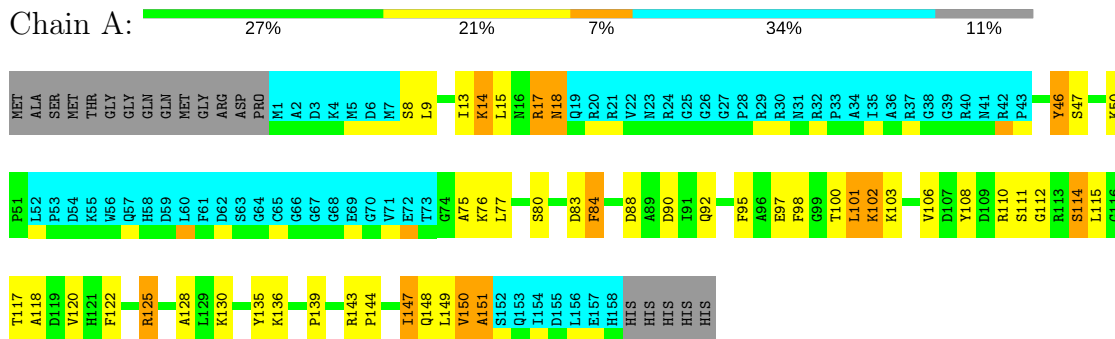
### 4.2.4 Score per residue for model 4

- Molecule 1: RNA and export factor binding protein 2



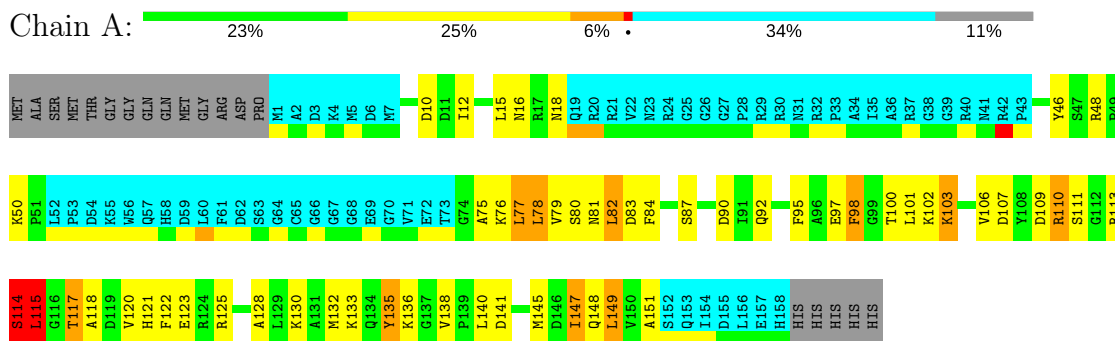
### 4.2.5 Score per residue for model 5

- Molecule 1: RNA and export factor binding protein 2



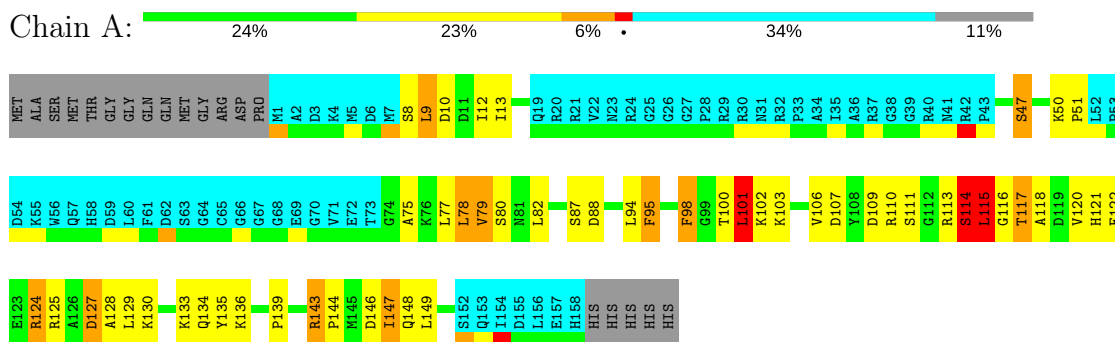
### 4.2.6 Score per residue for model 6

- Molecule 1: RNA and export factor binding protein 2



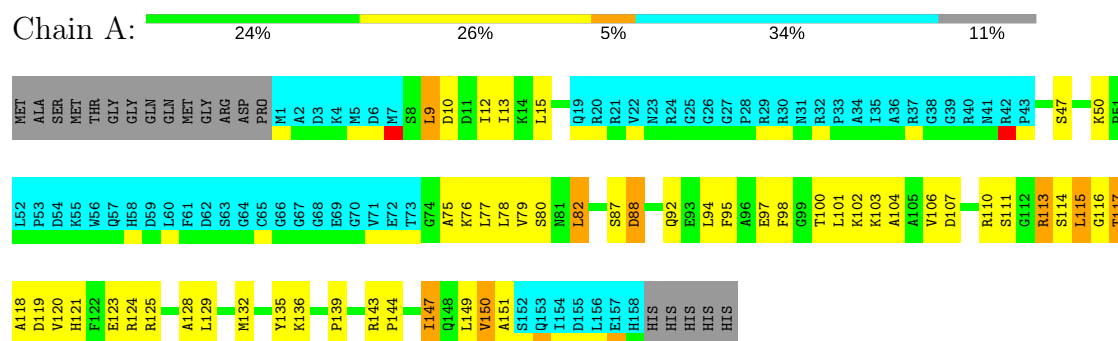
### 4.2.7 Score per residue for model 7 (medoid)

- Molecule 1: RNA and export factor binding protein 2



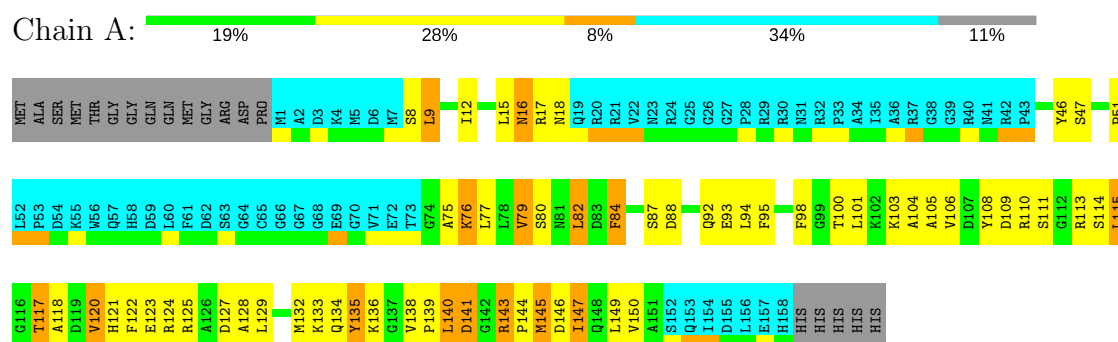
### 4.2.8 Score per residue for model 8

- Molecule 1: RNA and export factor binding protein 2



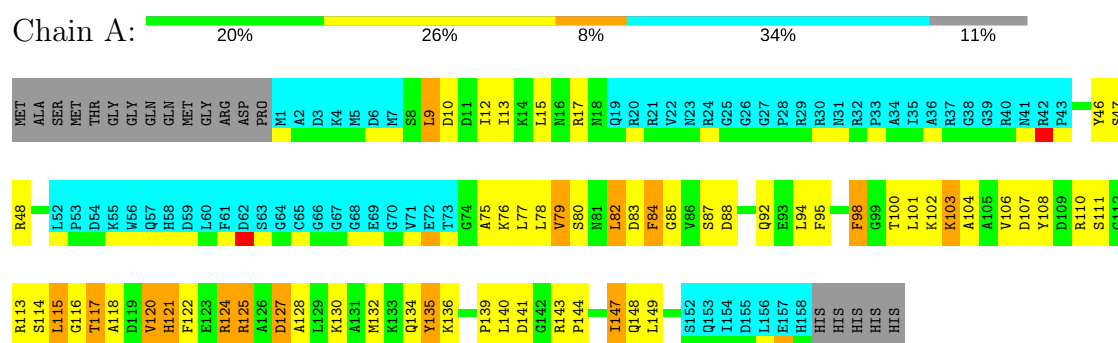
### 4.2.9 Score per residue for model 9

- Molecule 1: RNA and export factor binding protein 2



### 4.2.10 Score per residue for model 10

- Molecule 1: RNA and export factor binding protein 2

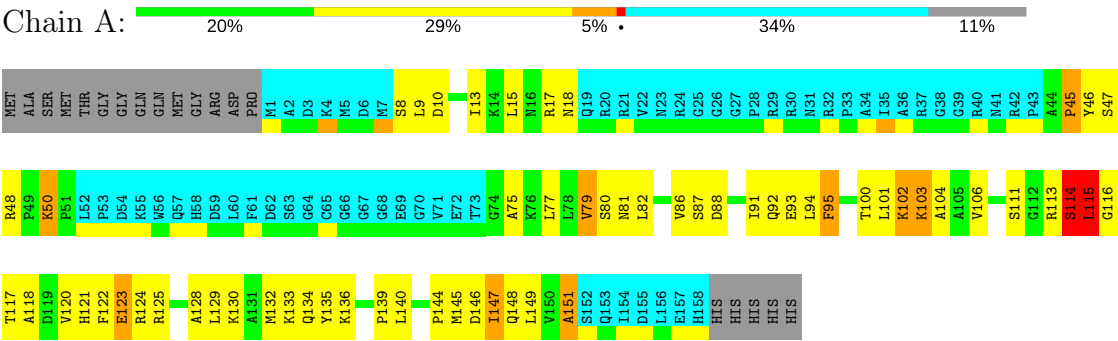






4.2.14 Score per residue for model 14

- Molecule 1: RNA and export factor binding protein 2



## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *ARIA protocol (Nilges et al., 1997, J. Mol. Biol. 269, 408-422) used for structure calculation..*

Of the 20 calculated structures, 14 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	structure solution	1.1
CNS	structure solution	1.0
CNS	refinement	1.0

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	751	762	761	29±6
All	All	10514	10668	10654	410

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 19.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:128:ALA:HA	1:A:149:LEU:HD22	0.90	1.43	2	4
1:A:128:ALA:HB1	1:A:149:LEU:HD11	0.89	1.44	1	1
1:A:139:PRO:HA	1:A:144:PRO:HA	0.89	1.44	4	12
1:A:136:LYS:HA	1:A:147:ILE:HG13	0.88	1.45	14	5
1:A:77:LEU:HB3	1:A:120:VAL:HG22	0.86	1.47	9	13
1:A:101:LEU:HD11	1:A:120:VAL:HB	0.84	1.46	7	6
1:A:77:LEU:HD23	1:A:149:LEU:HG	0.84	1.47	2	4
1:A:85:GLY:HA2	1:A:108:TYR:HA	0.83	1.50	4	1
1:A:95:PHE:HB3	1:A:101:LEU:HG	0.82	1.46	13	4
1:A:86:VAL:HG23	1:A:116:GLY:HA2	0.78	1.53	4	1
1:A:107:ASP:HB3	1:A:117:THR:HB	0.76	1.54	8	3
1:A:75:ALA:HB1	1:A:128:ALA:HB2	0.74	1.59	6	8
1:A:76:LYS:HE3	1:A:78:LEU:HD11	0.74	1.59	10	1
1:A:143:ARG:HG3	1:A:144:PRO:HD2	0.74	1.58	7	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:79:VAL:O	1:A:117:THR:HA	0.73	1.83	1	12
1:A:77:LEU:HD21	1:A:147:ILE:HG22	0.71	1.62	7	3
1:A:109:ASP:HB3	1:A:115:LEU:HD21	0.71	1.61	1	1
1:A:45:PRO:HD2	1:A:49:PRO:HA	0.71	1.61	3	3
1:A:77:LEU:HB3	1:A:120:VAL:HB	0.70	1.61	1	1
1:A:80:SER:HB2	1:A:146:ASP:HB3	0.70	1.63	13	3
1:A:124:ARG:HG3	1:A:127:ASP:HB2	0.70	1.64	10	1
1:A:92:GLN:HA	1:A:101:LEU:HD23	0.70	1.63	6	3
1:A:103:LYS:HG3	1:A:121:HIS:HB2	0.69	1.63	3	4
1:A:95:PHE:HA	1:A:98:PHE:HB2	0.69	1.65	8	9
1:A:77:LEU:HB3	1:A:120:VAL:CG2	0.69	2.18	13	2
1:A:101:LEU:HD11	1:A:120:VAL:HG13	0.68	1.64	1	1
1:A:145:MET:HG3	1:A:147:ILE:HG13	0.68	1.66	11	2
1:A:77:LEU:HD12	1:A:120:VAL:HG13	0.68	1.64	10	5
1:A:77:LEU:HD13	1:A:78:LEU:N	0.67	2.04	3	1
1:A:95:PHE:HB2	1:A:101:LEU:HD22	0.67	1.64	6	3
1:A:139:PRO:HA	1:A:144:PRO:CA	0.67	2.18	7	7
1:A:76:LYS:HB3	1:A:150:VAL:HG13	0.66	1.65	8	2
1:A:87:SER:HA	1:A:108:TYR:HE1	0.66	1.50	4	1
1:A:79:VAL:HG23	1:A:118:ALA:HB3	0.66	1.67	9	4
1:A:79:VAL:HG22	1:A:118:ALA:HB3	0.66	1.66	13	1
1:A:85:GLY:H	1:A:117:THR:HB	0.66	1.50	4	1
1:A:82:LEU:HB2	1:A:116:GLY:HA3	0.66	1.66	12	3
1:A:103:LYS:HG2	1:A:121:HIS:HB2	0.65	1.68	11	2
1:A:77:LEU:HD21	1:A:149:LEU:HD23	0.64	1.69	13	1
1:A:109:ASP:HB2	1:A:115:LEU:HD13	0.64	1.68	4	3
1:A:109:ASP:HB3	1:A:113:ARG:H	0.64	1.53	11	1
1:A:77:LEU:HD12	1:A:120:VAL:HG22	0.63	1.70	11	7
1:A:106:VAL:HG12	1:A:118:ALA:HB2	0.62	1.71	12	4
1:A:77:LEU:HD11	1:A:79:VAL:HG13	0.62	1.70	11	4
1:A:79:VAL:HG12	1:A:147:ILE:HG12	0.62	1.71	11	1
1:A:84:PHE:HA	1:A:117:THR:HB	0.62	1.71	4	1
1:A:85:GLY:HA2	1:A:108:TYR:CA	0.61	2.25	4	1
1:A:85:GLY:HA2	1:A:107:ASP:O	0.61	1.95	4	1
1:A:88:ASP:HA	1:A:106:VAL:HG23	0.61	1.73	14	2
1:A:79:VAL:HG12	1:A:147:ILE:HG23	0.59	1.74	6	2
1:A:10:ASP:HA	1:A:13:ILE:HD12	0.59	1.75	7	8
1:A:95:PHE:CB	1:A:101:LEU:HD22	0.59	2.27	6	4
1:A:114:SER:C	1:A:115:LEU:HG	0.59	2.17	7	2
1:A:95:PHE:HB3	1:A:101:LEU:HD11	0.58	1.74	14	1
1:A:85:GLY:N	1:A:117:THR:HB	0.57	2.14	4	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:85:GLY:CA	1:A:108:TYR:HA	0.57	2.28	4	1
1:A:108:TYR:HB3	1:A:112:GLY:HA2	0.57	1.76	5	1
1:A:96:ALA:HA	1:A:99:GLY:O	0.57	2.00	1	2
1:A:132:MET:HA	1:A:149:LEU:HD13	0.57	1.77	13	1
1:A:136:LYS:HA	1:A:147:ILE:HD13	0.57	1.76	4	2
1:A:13:ILE:O	1:A:17:ARG:HD2	0.56	2.00	5	1
1:A:78:LEU:HB2	1:A:148:GLN:HG2	0.56	1.76	11	1
1:A:95:PHE:HB3	1:A:101:LEU:HD22	0.56	1.77	7	2
1:A:106:VAL:HA	1:A:118:ALA:HA	0.56	1.78	11	5
1:A:75:ALA:HA	1:A:151:ALA:HB2	0.56	1.78	14	1
1:A:86:VAL:N	1:A:117:THR:H	0.55	1.99	4	1
1:A:9:LEU:HA	1:A:12:ILE:HD12	0.55	1.77	10	6
1:A:135:TYR:O	1:A:138:VAL:HG23	0.55	2.01	6	1
1:A:76:LYS:HD2	1:A:150:VAL:HB	0.55	1.78	9	1
1:A:128:ALA:O	1:A:149:LEU:HG	0.55	2.02	8	3
1:A:135:TYR:O	1:A:138:VAL:HB	0.55	2.02	4	1
1:A:88:ASP:HB3	1:A:106:VAL:HG21	0.55	1.78	8	1
1:A:76:LYS:HG3	1:A:150:VAL:O	0.54	2.02	2	1
1:A:132:MET:HB2	1:A:149:LEU:HD13	0.54	1.79	13	1
1:A:145:MET:HG3	1:A:147:ILE:CD1	0.54	2.32	1	2
1:A:78:LEU:O	1:A:147:ILE:HA	0.54	2.03	1	2
1:A:140:LEU:HB2	1:A:145:MET:HB3	0.54	1.80	9	2
1:A:136:LYS:HD3	1:A:147:ILE:HG12	0.54	1.80	7	1
1:A:78:LEU:HB2	1:A:148:GLN:HB2	0.54	1.78	13	2
1:A:132:MET:CB	1:A:149:LEU:HD13	0.54	2.33	13	1
1:A:92:GLN:HA	1:A:101:LEU:CD1	0.54	2.33	13	1
1:A:91:ILE:HA	1:A:94:LEU:HD12	0.53	1.78	14	1
1:A:128:ALA:CA	1:A:149:LEU:HD22	0.53	2.28	2	2
1:A:138:VAL:O	1:A:145:MET:HG2	0.53	2.03	3	1
1:A:76:LYS:HE2	1:A:150:VAL:HG23	0.53	1.80	2	1
1:A:87:SER:HA	1:A:108:TYR:CE1	0.53	2.37	4	1
1:A:124:ARG:HB2	1:A:127:ASP:HB2	0.53	1.81	7	2
1:A:80:SER:O	1:A:81:ASN:HB2	0.52	2.04	1	2
1:A:88:ASP:HA	1:A:106:VAL:HB	0.52	1.81	2	1
1:A:85:GLY:HA3	1:A:115:LEU:HD22	0.52	1.81	4	1
1:A:86:VAL:HB	1:A:117:THR:N	0.52	2.20	4	1
1:A:79:VAL:HG13	1:A:118:ALA:HB3	0.52	1.82	8	1
1:A:79:VAL:CG1	1:A:118:ALA:HB3	0.52	2.35	8	1
1:A:101:LEU:HD11	1:A:104:ALA:HB2	0.52	1.81	3	1
1:A:84:PHE:CA	1:A:117:THR:HB	0.52	2.35	4	1
1:A:76:LYS:HG3	1:A:119:ASP:OD1	0.51	2.05	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:86:VAL:HG12	1:A:118:ALA:HB2	0.51	1.82	4	1
1:A:104:ALA:HA	1:A:119:ASP:O	0.51	2.05	1	1
1:A:136:LYS:HA	1:A:147:ILE:CG1	0.51	2.28	14	4
1:A:101:LEU:HD21	1:A:104:ALA:HB2	0.51	1.82	9	2
1:A:114:SER:O	1:A:115:LEU:HG	0.51	2.05	7	2
1:A:8:SER:HB2	1:A:11:ASP:HB2	0.51	1.81	11	1
1:A:75:ALA:O	1:A:121:HIS:HA	0.50	2.06	12	5
1:A:132:MET:HG2	1:A:133:LYS:HD3	0.50	1.83	13	1
1:A:115:LEU:HD21	1:A:117:THR:HB	0.50	1.83	14	1
1:A:9:LEU:HG	1:A:12:ILE:HD12	0.50	1.82	1	1
1:A:82:LEU:HG	1:A:85:GLY:HA2	0.50	1.84	11	1
1:A:128:ALA:HB1	1:A:149:LEU:HD12	0.49	1.83	9	2
1:A:77:LEU:HD12	1:A:120:VAL:CG1	0.49	2.37	14	4
1:A:114:SER:O	1:A:116:GLY:N	0.49	2.44	14	2
1:A:95:PHE:HA	1:A:98:PHE:HB3	0.49	1.85	9	1
1:A:77:LEU:O	1:A:119:ASP:HA	0.49	2.08	4	1
1:A:136:LYS:CA	1:A:147:ILE:HG13	0.49	2.38	7	2
1:A:12:ILE:HG22	1:A:16:ASN:ND2	0.49	2.22	6	2
1:A:82:LEU:HA	1:A:116:GLY:HA3	0.49	1.84	10	1
1:A:131:ALA:HA	1:A:135:TYR:CD2	0.48	2.43	3	1
1:A:76:LYS:HB3	1:A:150:VAL:O	0.48	2.09	1	1
1:A:78:LEU:HD11	1:A:148:GLN:HB2	0.48	1.84	4	1
1:A:139:PRO:CA	1:A:144:PRO:HA	0.48	2.35	3	1
1:A:125:ARG:HB2	1:A:151:ALA:HB1	0.48	1.86	5	1
1:A:77:LEU:HD23	1:A:149:LEU:CG	0.48	2.28	2	2
1:A:77:LEU:HD13	1:A:77:LEU:C	0.48	2.29	4	3
1:A:115:LEU:HD13	1:A:117:THR:OG1	0.48	2.09	8	2
1:A:132:MET:CA	1:A:149:LEU:HD13	0.48	2.38	13	1
1:A:145:MET:HG3	1:A:147:ILE:HD11	0.47	1.86	1	1
1:A:81:ASN:O	1:A:82:LEU:HB2	0.47	2.08	6	1
1:A:82:LEU:CB	1:A:116:GLY:HA3	0.47	2.40	1	1
1:A:77:LEU:CD1	1:A:79:VAL:HG22	0.47	2.39	14	2
1:A:44:ALA:HA	1:A:49:PRO:HG3	0.47	1.86	12	1
1:A:103:LYS:HG3	1:A:121:HIS:CB	0.47	2.40	6	3
1:A:104:ALA:HA	1:A:120:VAL:HA	0.47	1.84	3	1
1:A:76:LYS:HE2	1:A:78:LEU:HD21	0.47	1.86	6	1
1:A:130:LYS:O	1:A:134:GLN:HB2	0.47	2.09	1	2
1:A:101:LEU:HD21	1:A:120:VAL:HB	0.47	1.85	3	1
1:A:77:LEU:C	1:A:77:LEU:HD13	0.46	2.30	8	5
1:A:95:PHE:HA	1:A:98:PHE:CB	0.46	2.40	9	1
1:A:86:VAL:O	1:A:106:VAL:HB	0.46	2.10	4	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:77:LEU:CB	1:A:120:VAL:HG22	0.46	2.33	10	1
1:A:101:LEU:HD13	1:A:104:ALA:CB	0.46	2.40	14	1
1:A:76:LYS:HE3	1:A:78:LEU:HD21	0.46	1.87	8	1
1:A:108:TYR:HA	1:A:115:LEU:HD11	0.46	1.86	11	1
1:A:82:LEU:HB3	1:A:86:VAL:HG21	0.46	1.88	1	1
1:A:88:ASP:HB2	1:A:105:ALA:HA	0.46	1.86	9	1
1:A:132:MET:O	1:A:136:LYS:HB3	0.46	2.09	6	2
1:A:108:TYR:HD2	1:A:113:ARG:HD2	0.46	1.70	1	1
1:A:102:LYS:HB3	1:A:121:HIS:O	0.46	2.10	11	1
1:A:104:ALA:HB2	1:A:120:VAL:HB	0.46	1.88	14	2
1:A:77:LEU:HD22	1:A:78:LEU:H	0.46	1.71	3	1
1:A:132:MET:HB3	1:A:149:LEU:HB2	0.46	1.88	14	2
1:A:85:GLY:HA2	1:A:107:ASP:C	0.45	2.32	4	1
1:A:15:LEU:HA	1:A:18:ASN:OD1	0.45	2.11	14	1
1:A:132:MET:O	1:A:136:LYS:HB2	0.45	2.11	1	2
1:A:109:ASP:HB2	1:A:115:LEU:HD21	0.45	1.87	11	1
1:A:122:PHE:N	1:A:122:PHE:CD1	0.45	2.85	1	1
1:A:92:GLN:HG3	1:A:104:ALA:HB3	0.45	1.87	10	1
1:A:76:LYS:HB3	1:A:121:HIS:CE1	0.45	2.47	2	1
1:A:77:LEU:HD11	1:A:147:ILE:HG22	0.44	1.89	1	1
1:A:75:ALA:HB3	1:A:122:PHE:HD1	0.44	1.72	2	1
1:A:132:MET:HG3	1:A:133:LYS:HD2	0.44	1.89	3	1
1:A:84:PHE:HE1	1:A:141:ASP:HB2	0.44	1.71	9	1
1:A:108:TYR:HA	1:A:113:ARG:O	0.44	2.11	10	1
1:A:79:VAL:CG2	1:A:118:ALA:HB3	0.44	2.42	9	2
1:A:79:VAL:HA	1:A:146:ASP:O	0.44	2.12	14	2
1:A:132:MET:HA	1:A:136:LYS:CB	0.44	2.43	14	1
1:A:92:GLN:HA	1:A:101:LEU:HD12	0.44	1.90	13	1
1:A:88:ASP:HA	1:A:106:VAL:HG13	0.43	1.88	5	2
1:A:89:ALA:O	1:A:93:GLU:HB2	0.43	2.14	12	1
1:A:92:GLN:HG3	1:A:101:LEU:HD23	0.43	1.89	1	1
1:A:82:LEU:HD22	1:A:91:ILE:HD11	0.43	1.90	3	1
1:A:101:LEU:HD13	1:A:104:ALA:HB3	0.43	1.90	14	1
1:A:136:LYS:HG3	1:A:147:ILE:O	0.43	2.13	10	2
1:A:82:LEU:HG	1:A:116:GLY:HA3	0.43	1.90	8	1
1:A:80:SER:HB2	1:A:146:ASP:CB	0.43	2.43	9	1
1:A:129:LEU:HG	1:A:133:LYS:HE2	0.43	1.90	2	1
1:A:86:VAL:HB	1:A:117:THR:H	0.42	1.74	4	1
1:A:109:ASP:CB	1:A:115:LEU:HD21	0.42	2.41	1	1
1:A:132:MET:HA	1:A:136:LYS:HB3	0.42	1.90	10	1
1:A:79:VAL:HG23	1:A:118:ALA:H	0.42	1.74	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:102:LYS:HG3	1:A:123:GLU:HG3	0.42	1.91	13	1
1:A:78:LEU:HB2	1:A:148:GLN:CB	0.42	2.44	3	1
1:A:95:PHE:HB3	1:A:101:LEU:CG	0.42	2.32	13	2
1:A:125:ARG:HB3	1:A:151:ALA:HB1	0.42	1.90	4	1
1:A:77:LEU:HG	1:A:95:PHE:CE2	0.42	2.49	4	1
1:A:109:ASP:OD2	1:A:110:ARG:HD2	0.42	2.14	1	1
1:A:77:LEU:HD22	1:A:148:GLN:O	0.42	2.14	3	1
1:A:105:ALA:O	1:A:118:ALA:HB1	0.42	2.15	3	1
1:A:48:ARG:HG3	1:A:49:PRO:HD2	0.42	1.91	2	1
1:A:77:LEU:HG	1:A:95:PHE:CZ	0.42	2.49	13	3
1:A:78:LEU:HA	1:A:118:ALA:O	0.42	2.14	13	2
1:A:88:ASP:HA	1:A:106:VAL:CG2	0.42	2.45	14	1
1:A:103:LYS:CG	1:A:121:HIS:HB2	0.42	2.44	12	1
1:A:77:LEU:HD23	1:A:149:LEU:HD22	0.41	1.91	14	1
1:A:101:LEU:HA	1:A:101:LEU:HD23	0.41	1.69	5	3
1:A:149:LEU:HD13	1:A:149:LEU:HA	0.41	1.78	3	1
1:A:109:ASP:HB3	1:A:113:ARG:N	0.41	2.25	11	1
1:A:101:LEU:HD11	1:A:120:VAL:CG1	0.41	2.42	1	1
1:A:132:MET:HB3	1:A:149:LEU:HD13	0.41	1.91	1	1
1:A:102:LYS:HD3	1:A:102:LYS:HA	0.41	1.54	2	1
1:A:10:ASP:O	1:A:14:LYS:HG2	0.41	2.15	4	4
1:A:128:ALA:HB1	1:A:149:LEU:HB3	0.41	1.93	4	1
1:A:87:SER:N	1:A:90:ASP:HB2	0.41	2.30	13	1
1:A:77:LEU:HA	1:A:149:LEU:HD13	0.41	1.91	3	1
1:A:102:LYS:HB2	1:A:121:HIS:O	0.41	2.16	1	1
1:A:14:LYS:O	1:A:18:ASN:HB3	0.41	2.15	5	1
1:A:98:PHE:HE2	1:A:135:TYR:CD2	0.41	2.33	10	1
1:A:143:ARG:HD2	1:A:144:PRO:O	0.41	2.16	10	1
1:A:88:ASP:CB	1:A:106:VAL:HG21	0.41	2.45	8	1
1:A:109:ASP:OD1	1:A:110:ARG:HD2	0.41	2.16	6	1
1:A:115:LEU:N	1:A:115:LEU:HD12	0.40	2.32	1	1
1:A:131:ALA:HA	1:A:135:TYR:CE2	0.40	2.51	1	1
1:A:85:GLY:O	1:A:115:LEU:N	0.40	2.55	4	1
1:A:82:LEU:HA	1:A:115:LEU:HA	0.40	1.94	13	1
1:A:135:TYR:O	1:A:138:VAL:HG12	0.40	2.17	9	1
1:A:106:VAL:HG13	1:A:117:THR:O	0.40	2.15	10	1
1:A:102:LYS:HG3	1:A:123:GLU:HG2	0.40	1.92	14	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	97/177 (55%)	80±3 (83±3%)	13±2 (13±3%)	4±2 (4±2%)	6	32
All	All	1358/2478 (55%)	1125 (83%)	178 (13%)	55 (4%)	6	32

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	115	LEU	7
1	A	141	ASP	6
1	A	114	SER	6
1	A	113	ARG	5
1	A	81	ASN	4
1	A	151	ALA	4
1	A	51	PRO	3
1	A	82	LEU	3
1	A	45	PRO	2
1	A	85	GLY	2
1	A	83	ASP	2
1	A	46	TYR	2
1	A	101	LEU	2
1	A	84	PHE	2
1	A	117	THR	1
1	A	142	GLY	1
1	A	147	ILE	1
1	A	79	VAL	1
1	A	112	GLY	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	79/142 (56%)	46±4 (58±5%)	33±4 (42±5%)	0 4
All	All	1106/1988 (56%)	638 (58%)	468 (42%)	0 4

All 69 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	103	LYS	14
1	A	100	THR	14
1	A	122	PHE	13
1	A	87	SER	13
1	A	147	ILE	13
1	A	111	SER	12
1	A	115	LEU	12
1	A	50	LYS	11
1	A	47	SER	11
1	A	102	LYS	11
1	A	8	SER	11
1	A	80	SER	10
1	A	135	TYR	10
1	A	123	GLU	10
1	A	9	LEU	10
1	A	82	LEU	10
1	A	130	LYS	10
1	A	114	SER	10
1	A	97	GLU	9
1	A	133	LYS	9
1	A	46	TYR	9
1	A	117	THR	9
1	A	124	ARG	9
1	A	143	ARG	8
1	A	125	ARG	8
1	A	15	LEU	8
1	A	134	GLN	8
1	A	110	ARG	8
1	A	140	LEU	8
1	A	48	ARG	7
1	A	113	ARG	7
1	A	94	LEU	7
1	A	148	GLN	7
1	A	84	PHE	7
1	A	127	ASP	7
1	A	95	PHE	6

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Mol	Chain	Res	Type	Models (Total)
1	A	79	VAL	6
1	A	90	ASP	6
1	A	88	ASP	6
1	A	101	LEU	6
1	A	121	HIS	5
1	A	93	GLU	5
1	A	145	MET	5
1	A	17	ARG	5
1	A	107	ASP	5
1	A	129	LEU	5
1	A	98	PHE	4
1	A	120	VAL	4
1	A	92	GLN	4
1	A	78	LEU	4
1	A	77	LEU	4
1	A	149	LEU	4
1	A	119	ASP	4
1	A	150	VAL	4
1	A	83	ASP	4
1	A	136	LYS	4
1	A	86	VAL	4
1	A	109	ASP	4
1	A	10	ASP	3
1	A	16	ASN	3
1	A	132	MET	2
1	A	76	LYS	2
1	A	18	ASN	2
1	A	11	ASP	2
1	A	141	ASP	2
1	A	81	ASN	1
1	A	108	TYR	1
1	A	138	VAL	1
1	A	14	LYS	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