Full wwPDB NMR Structure Validation Report (i)

Feb 24, 2024 – 08:33 AM EST

PDB ID	:	1XPN
Title	:	NMR structure of P. aeruginosa protein PA1324: Northeast Structural Genomics Consortium target PaP1
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Deposited on	:	2004-10-08

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
https://www.wwpdb.org/validation/2017/NMRValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (i)) were used in the production of this report:

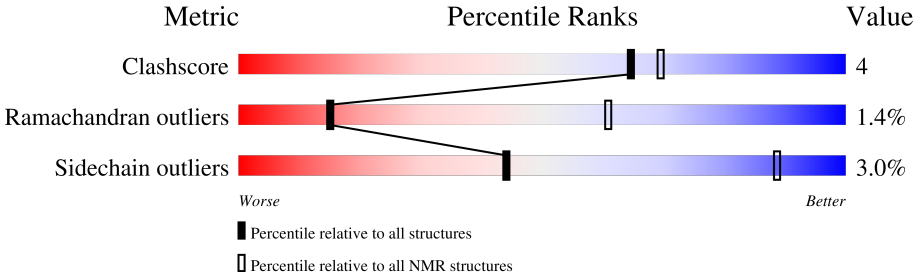
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 >=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 <=5%

Mol	Chain	Length	Quality of chain
1	A	170	84% • 15%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 11 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy, fewest violations, closest to the average.*

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:27-A:170 (144)	1.18	11

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 4 clusters and 3 single-model clusters were found.

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

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2546 atoms, of which 1248 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called hypothetical protein PA1324.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	170	2546	804	1248	233	259	2	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP Q9I420
A	2	SER	-	expression tag	UNP Q9I420
A	3	SER	-	expression tag	UNP Q9I420
A	4	HIS	-	expression tag	UNP Q9I420
A	5	HIS	-	expression tag	UNP Q9I420
A	6	HIS	-	expression tag	UNP Q9I420
A	7	HIS	-	expression tag	UNP Q9I420
A	8	HIS	-	expression tag	UNP Q9I420
A	9	HIS	-	expression tag	UNP Q9I420
A	10	SER	-	expression tag	UNP Q9I420
A	11	SER	-	expression tag	UNP Q9I420
A	12	GLY	-	expression tag	UNP Q9I420
A	13	LEU	-	expression tag	UNP Q9I420
A	14	VAL	-	expression tag	UNP Q9I420
A	15	PRO	-	expression tag	UNP Q9I420
A	16	ARG	-	expression tag	UNP Q9I420
A	17	GLY	-	expression tag	UNP Q9I420
A	18	SER	-	expression tag	UNP Q9I420
A	19	HIS	-	expression tag	UNP Q9I420
A	20	MET	-	expression tag	UNP Q9I420
A	43	GLY	GLU	conflict	UNP Q9I420

4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide 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: hypothetical protein PA1324

Chain A: 84% • 15%

G1 S2 S3 H4 H5 H6 H7 H8 H9 S10 S11 G12 L13 V14 P15 R16 G17 S18 H19 M20 A21 S22 N23 P24 N25 D26 G48 K61 R170

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: hypothetical protein PA1324

Chain A: 71% 14% 15%

G1 S2 S3 H4 H5 H6 H7 H8 H9 S10 S11 G12 L13 V14 P15 R16 G17 S18 H19 M20 A21 S22 N23 P24 N25 D26 L27 P28 Q40 V41 G42 Y51 L52 T53 S56 K61 G62 T63 Y75 M76 N85 V99 L100 R104 V108 D114 H123 G143

T151 K154 P163 L168 T169 R170

4.2.2 Score per residue for model 2

• Molecule 1: hypothetical protein PA1324

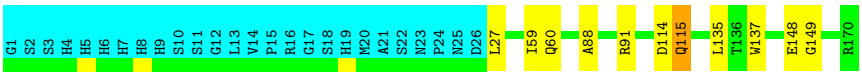
Chain A: 69% 15% • 15%

G1 S2 S3 H4 H5 H6 H7 H8 H9 S10 S11 G12 L13 V14 P15 R16 G17 S18 H19 M20 A21 S22 N23 P24 N25 D26 Q39 Q40 V41 G42 G48 D49 L50 K61 E70 T73 M76 N85 L86 D87 D101 R104 V108 D114 Q115 H123 N126

4.2.3 Score per residue for model 3

• Molecule 1: hypothetical protein PA1324

Chain A:



4.2.4 Score per residue for model 4

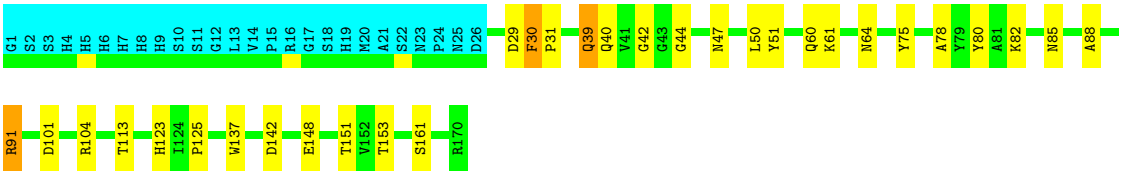
• Molecule 1: hypothetical protein PA1324

Chain A:

4.2.5 Score per residue for model 5

• Molecule 1: hypothetical protein PA1324

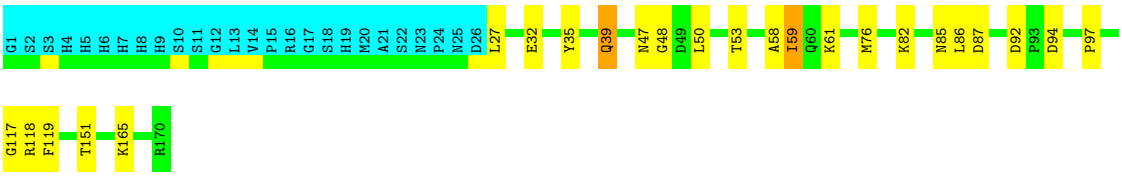
Chain A:



4.2.6 Score per residue for model 6

• Molecule 1: hypothetical protein PA1324

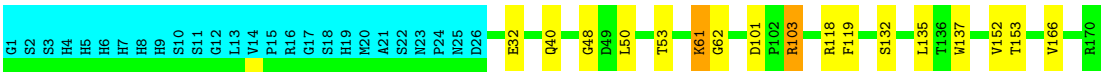
Chain A:



4.2.7 Score per residue for model 7

- Molecule 1: hypothetical protein PA1324

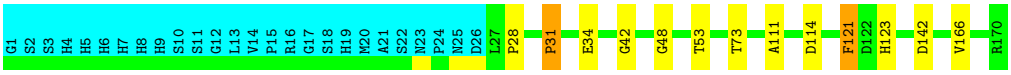
Chain A: 75% 9% • 15%



4.2.8 Score per residue for model 8

- Molecule 1: hypothetical protein PA1324

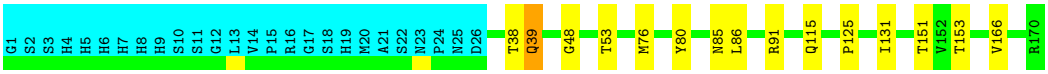
Chain A: 77% 6% • 15%



4.2.9 Score per residue for model 9

- Molecule 1: hypothetical protein PA1324

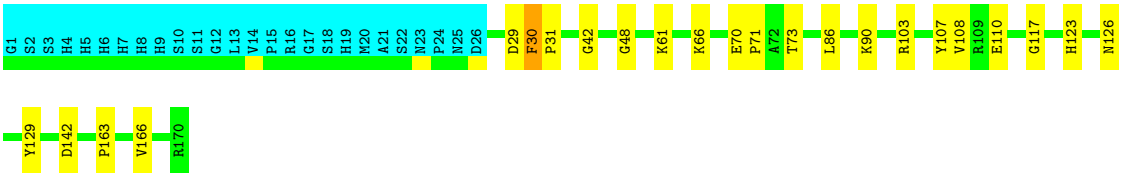
Chain A: 76% 8% • 15%



4.2.10 Score per residue for model 10

- Molecule 1: hypothetical protein PA1324

Chain A: 71% 13% • 15%



4.2.11 Score per residue for model 11 (medoid)

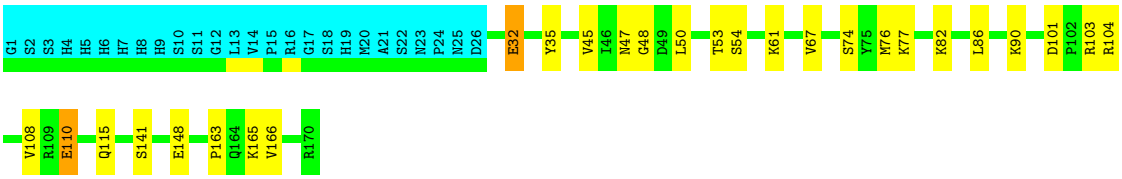
- Molecule 1: hypothetical protein PA1324

Chain A: 78% 6% 15%

4.2.12 Score per residue for model 12

- Molecule 1: hypothetical protein PA1324

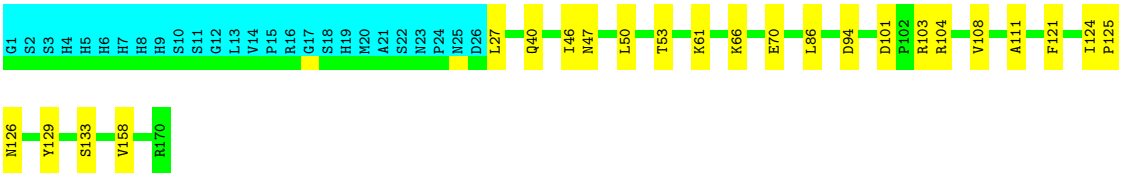
Chain A:



4.2.13 Score per residue for model 13

- Molecule 1: hypothetical protein PA1324

Chain A:



4.2.14 Score per residue for model 14

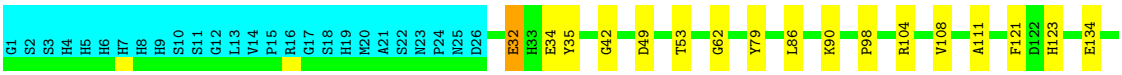
- Molecule 1: hypothetical protein PA1324

Chain A:

4.2.15 Score per residue for model 15

- Molecule 1: hypothetical protein PA1324

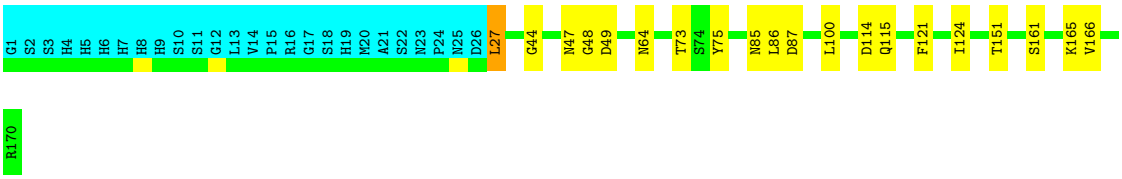
Chain A:



4.2.16 Score per residue for model 16

• Molecule 1: hypothetical protein PA1324

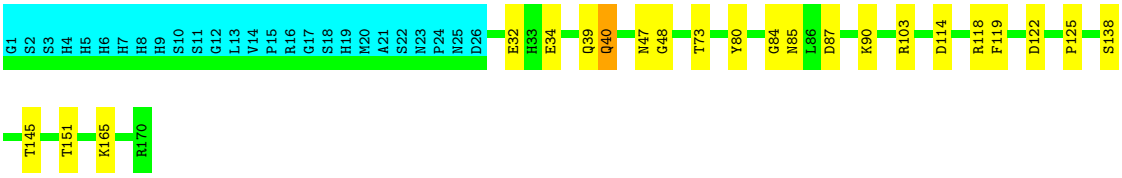
Chain A:



4.2.17 Score per residue for model 17

• Molecule 1: hypothetical protein PA1324

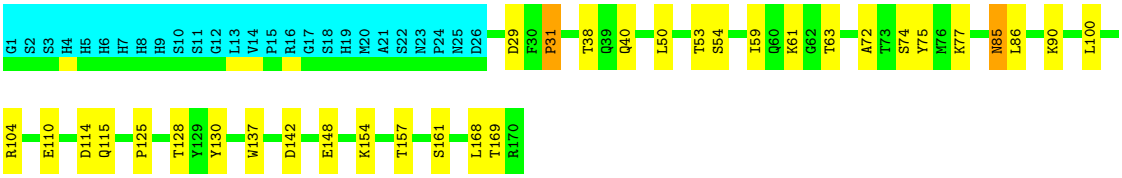
Chain A:



4.2.18 Score per residue for model 18

• Molecule 1: hypothetical protein PA1324

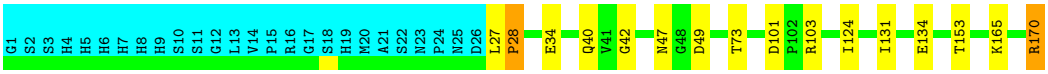
Chain A:



4.2.19 Score per residue for model 19

• Molecule 1: hypothetical protein PA1324

Chain A:



4.2.20 Score per residue for model 20

- Molecule 1: hypothetical protein PA1324

Chain A:

5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *Molecular Dynamics and Simulated Annealing*.

Of the 20 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy,structures with the fewest restraint violations*.

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

Software name	Classification	Version
NIH-Xplor	refinement	2.0.4
CNS	refinement	1.1
AutoStructure	structure solution	2.1.0

No chemical shift data was provided.

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	1103	1074	1072	9±4
All	All	22060	21480	21440	182

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 4.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:100:LEU:HB2	1:A:104:ARG:HE	0.76	1.39	18	1
1:A:39:GLN:HE21	1:A:39:GLN:HA	0.73	1.44	6	1
1:A:82:LYS:HE2	1:A:82:LYS:HA	0.65	1.66	12	1
1:A:76:MET:HG3	1:A:104:ARG:HH21	0.65	1.51	2	1
1:A:86:LEU:HA	1:A:90:LYS:HE3	0.65	1.67	4	1
1:A:86:LEU:HA	1:A:90:LYS:HE2	0.60	1.72	18	2
1:A:44:GLY:HA2	1:A:161:SER:HA	0.60	1.71	5	2
1:A:104:ARG:O	1:A:108:VAL:HG23	0.60	1.95	12	6
1:A:70:GLU:HG2	1:A:108:VAL:HG22	0.57	1.74	10	1
1:A:46:ILE:HD11	1:A:158:VAL:HG21	0.57	1.75	13	1
1:A:85:ASN:HA	1:A:151:THR:O	0.56	2.01	6	7
1:A:34:GLU:OE1	1:A:73:THR:HB	0.56	2.00	17	4
1:A:67:VAL:O	1:A:110:GLU:HA	0.55	2.00	12	1
1:A:48:GLY:HA2	1:A:166:VAL:O	0.55	2.01	11	7

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:29:ASP:O	1:A:30:PHE:HB2	0.55	2.01	10	2
1:A:47:ASN:O	1:A:165:LYS:HA	0.55	2.01	16	5
1:A:80:TYR:HE2	1:A:153:THR:HG21	0.55	1.61	4	1
1:A:30:PHE:HB2	1:A:31:PRO:HD3	0.54	1.79	20	1
1:A:29:ASP:O	1:A:31:PRO:HD3	0.54	2.03	18	2
1:A:50:LEU:O	1:A:61:LYS:HD3	0.54	2.03	12	3
1:A:86:LEU:HD23	1:A:87:ASP:O	0.53	2.02	6	2
1:A:28:PRO:HB3	1:A:75:TYR:HA	0.53	1.79	1	1
1:A:78:ALA:O	1:A:82:LYS:HG2	0.53	2.04	5	1
1:A:29:ASP:C	1:A:31:PRO:HD2	0.52	2.25	10	1
1:A:42:GLY:O	1:A:123:HIS:HA	0.52	2.05	15	8
1:A:101:ASP:OD2	1:A:103:ARG:HB2	0.52	2.04	13	2
1:A:73:THR:HG23	1:A:75:TYR:H	0.52	1.64	16	1
1:A:58:ALA:O	1:A:59:ILE:HB	0.52	2.05	20	2
1:A:103:ARG:HE	1:A:103:ARG:HA	0.52	1.64	10	1
1:A:111:ALA:HB2	1:A:121:PHE:CZ	0.52	2.40	15	2
1:A:114:ASP:HB2	1:A:118:ARG:HB2	0.51	1.79	17	1
1:A:131:ILE:O	1:A:153:THR:HA	0.51	2.05	19	4
1:A:135:LEU:HG	1:A:137:TRP:CZ3	0.51	2.40	7	1
1:A:61:LYS:HB3	1:A:117:GLY:HA3	0.50	1.82	6	2
1:A:50:LEU:HG	1:A:61:LYS:O	0.50	2.06	6	5
1:A:80:TYR:CE2	1:A:153:THR:HG21	0.50	2.40	4	1
1:A:48:GLY:HA3	1:A:166:VAL:O	0.49	2.06	12	2
1:A:121:PHE:CB	1:A:124:ILE:HD11	0.49	2.38	16	1
1:A:82:LYS:HE2	1:A:97:PRO:HA	0.49	1.84	6	1
1:A:54:SER:HB2	1:A:148:GLU:OE2	0.49	2.08	12	1
1:A:135:LEU:O	1:A:149:GLY:HA3	0.49	2.07	3	1
1:A:28:PRO:HB2	1:A:31:PRO:HG2	0.49	1.84	8	1
1:A:154:LYS:HD2	1:A:168:LEU:HG	0.49	1.85	1	1
1:A:125:PRO:HB2	1:A:129:TYR:OH	0.48	2.08	13	1
1:A:39:GLN:HE21	1:A:39:GLN:CA	0.48	2.19	6	1
1:A:35:TYR:OH	1:A:76:MET:HG2	0.48	2.07	6	1
1:A:86:LEU:HG	1:A:90:LYS:HB3	0.48	1.84	15	1
1:A:80:TYR:HE1	1:A:153:THR:HG21	0.48	1.69	5	2
1:A:31:PRO:HD2	1:A:75:TYR:HB2	0.48	1.86	18	1
1:A:101:ASP:OD1	1:A:103:ARG:HG3	0.48	2.08	12	1
1:A:74:SER:O	1:A:77:LYS:HG2	0.47	2.09	12	1
1:A:100:LEU:HB2	1:A:104:ARG:NE	0.47	2.18	18	1
1:A:32:GLU:OE2	1:A:103:ARG:HD2	0.47	2.10	12	1
1:A:82:LYS:HA	1:A:82:LYS:CE	0.47	2.40	12	1
1:A:40:GLN:HE21	1:A:40:GLN:HA	0.46	1.71	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:70:GLU:HG3	1:A:130:TYR:HB2	0.46	1.88	2	1
1:A:137:TRP:NE1	1:A:148:GLU:HB2	0.46	2.26	18	3
1:A:137:TRP:CE2	1:A:148:GLU:HB2	0.46	2.45	3	3
1:A:60:GLN:HG3	1:A:137:TRP:CE2	0.46	2.46	3	1
1:A:66:LYS:O	1:A:133:SER:HA	0.46	2.11	13	1
1:A:75:TYR:OH	1:A:100:LEU:HA	0.46	2.10	1	1
1:A:48:GLY:HA3	1:A:119:PHE:CZ	0.45	2.47	7	3
1:A:39:GLN:O	1:A:40:GLN:HB2	0.45	2.11	5	2
1:A:51:TYR:HA	1:A:60:GLN:O	0.45	2.11	5	1
1:A:103:ARG:HD3	1:A:107:TYR:HE2	0.45	1.71	10	1
1:A:138:SER:OG	1:A:145:THR:HG22	0.45	2.12	2	1
1:A:64:ASN:HA	1:A:113:THR:O	0.45	2.11	5	2
1:A:86:LEU:HG	1:A:90:LYS:HG3	0.45	1.89	4	1
1:A:70:GLU:HB3	1:A:108:VAL:HG22	0.45	1.88	14	3
1:A:134:GLU:HA	1:A:150:GLY:O	0.45	2.12	15	1
1:A:86:LEU:HG	1:A:90:LYS:HB2	0.44	1.88	10	1
1:A:38:THR:O	1:A:125:PRO:HG2	0.44	2.12	18	1
1:A:72:ALA:HA	1:A:130:TYR:CE2	0.44	2.48	18	1
1:A:66:LYS:CE	1:A:110:GLU:HB3	0.44	2.43	10	1
1:A:40:GLN:O	1:A:125:PRO:HG3	0.44	2.13	5	1
1:A:79:TYR:CE1	1:A:98:PRO:HD2	0.44	2.47	15	1
1:A:114:ASP:CG	1:A:115:GLN:H	0.44	2.16	16	1
1:A:87:ASP:H	1:A:90:LYS:HD3	0.44	1.72	17	1
1:A:66:LYS:HE2	1:A:110:GLU:HB3	0.44	1.89	10	1
1:A:132:SER:HA	1:A:152:VAL:O	0.43	2.12	7	1
1:A:74:SER:O	1:A:77:LYS:HB2	0.43	2.13	18	1
1:A:27:LEU:HB3	1:A:28:PRO:CD	0.43	2.42	19	1
1:A:114:ASP:CG	1:A:115:GLN:N	0.43	2.72	3	4
1:A:86:LEU:HG	1:A:90:LYS:CB	0.43	2.44	15	1
1:A:101:ASP:O	1:A:104:ARG:HG2	0.43	2.13	5	1
1:A:35:TYR:HH	1:A:76:MET:HG2	0.43	1.74	6	1
1:A:80:TYR:O	1:A:84:GLY:HA3	0.43	2.14	17	1
1:A:32:GLU:HB2	1:A:35:TYR:CD2	0.43	2.48	15	1
1:A:138:SER:HB2	1:A:145:THR:HG22	0.43	1.91	17	1
1:A:31:PRO:HB2	1:A:75:TYR:HB2	0.42	1.92	5	1
1:A:135:LEU:HG	1:A:137:TRP:HZ3	0.42	1.73	7	1
1:A:134:GLU:HB3	1:A:151:THR:OG1	0.42	2.15	15	1
1:A:170:ARG:HG2	1:A:170:ARG:OXT	0.42	2.14	19	1
1:A:29:ASP:C	1:A:31:PRO:HD3	0.42	2.35	5	1
1:A:88:ALA:O	1:A:91:ARG:HB2	0.41	2.15	5	1
1:A:111:ALA:HB2	1:A:121:PHE:CE2	0.41	2.50	8	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:111:ALA:HB2	1:A:121:PHE:HZ	0.41	1.73	15	1
1:A:154:LYS:HE2	1:A:169:THR:O	0.41	2.15	18	1
1:A:51:TYR:HA	1:A:61:LYS:HA	0.41	1.92	1	1
1:A:35:TYR:CE2	1:A:104:ARG:HB3	0.41	2.51	12	1
1:A:82:LYS:NZ	1:A:82:LYS:HB3	0.41	2.30	14	1
1:A:39:GLN:O	1:A:125:PRO:HG2	0.41	2.16	17	1
1:A:85:ASN:HD22	1:A:85:ASN:H	0.41	1.58	18	1
1:A:30:PHE:N	1:A:31:PRO:CD	0.41	2.84	5	1
1:A:92:ASP:OD2	1:A:94:ASP:HB3	0.41	2.15	6	1
1:A:71:PRO:HA	1:A:129:TYR:CD2	0.41	2.51	10	1
1:A:134:GLU:HA	1:A:151:THR:HA	0.41	1.92	15	1
1:A:128:THR:HG22	1:A:157:THR:HB	0.41	1.93	18	1
1:A:101:ASP:O	1:A:104:ARG:HB2	0.40	2.16	2	1
1:A:39:GLN:HB2	1:A:125:PRO:CB	0.40	2.45	9	1
1:A:45:VAL:O	1:A:163:PRO:HA	0.40	2.16	12	1
1:A:121:PHE:HB2	1:A:124:ILE:HD11	0.40	1.92	16	1
1:A:101:ASP:OD1	1:A:103:ARG:HB2	0.40	2.16	7	1
1:A:80:TYR:CE1	1:A:153:THR:HG21	0.40	2.51	9	1
1:A:88:ALA:HA	1:A:91:ARG:HG2	0.40	1.92	3	1
1:A:140:GLN:HG2	1:A:145:THR:OG1	0.40	2.17	4	1
1:A:121:PHE:HB3	1:A:124:ILE:HD11	0.40	1.93	13	1
1:A:42:GLY:HA3	1:A:124:ILE:C	0.40	2.36	19	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	143/170 (84%)	130±3 (91±2%)	12±2 (8±2%)	2±1 (1±1%)	15	61
All	All	2860/3400 (84%)	2590 (91%)	230 (8%)	40 (1%)	15	61

All 16 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	142	ASP	6
1	A	59	ILE	5
1	A	62	GLY	4
1	A	27	LEU	4
1	A	40	GLN	3
1	A	114	ASP	2
1	A	163	PRO	2
1	A	115	GLN	2
1	A	30	PHE	2
1	A	31	PRO	2
1	A	32	GLU	2
1	A	28	PRO	2
1	A	143	GLY	1
1	A	61	LYS	1
1	A	117	GLY	1
1	A	141	SER	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	118/140 (84%)	114±2 (97±2%)	4±2 (3±2%)	44 89
All	All	2360/2800 (84%)	2289 (97%)	71 (3%)	44 89

All 37 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	53	THR	10
1	A	32	GLU	5
1	A	63	THR	3
1	A	76	MET	3
1	A	126	ASN	3
1	A	39	GLN	3
1	A	40	GLN	3
1	A	86	LEU	3
1	A	49	ASP	3
1	A	73	THR	2

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Mol	Chain	Res	Type	Models (Total)
1	A	47	ASN	2
1	A	91	ARG	2
1	A	118	ARG	2
1	A	103	ARG	2
1	A	110	GLU	2
1	A	100	LEU	2
1	A	56	SER	1
1	A	104	ARG	1
1	A	109	ARG	1
1	A	153	THR	1
1	A	121	PHE	1
1	A	38	THR	1
1	A	66	LYS	1
1	A	115	GLN	1
1	A	94	ASP	1
1	A	34	GLU	1
1	A	64	ASN	1
1	A	87	ASP	1
1	A	122	ASP	1
1	A	54	SER	1
1	A	85	ASN	1
1	A	161	SER	1
1	A	168	LEU	1
1	A	134	GLU	1
1	A	170	ARG	1
1	A	33	HIS	1
1	A	123	HIS	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 monosaccharides 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 (i)

No chemical shift data were provided