Full wwPDB NMR Structure Validation Report (i)

Feb 26, 2022 – 02:37 PM EST

PDB ID	:	2CSJ
Title	:	Solution structure of N-terminal PDZ domain from mouse TJP2
Authors	:	Suetake, T.; Hayashi, F.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on	:	2005-05-21

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/NMRValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

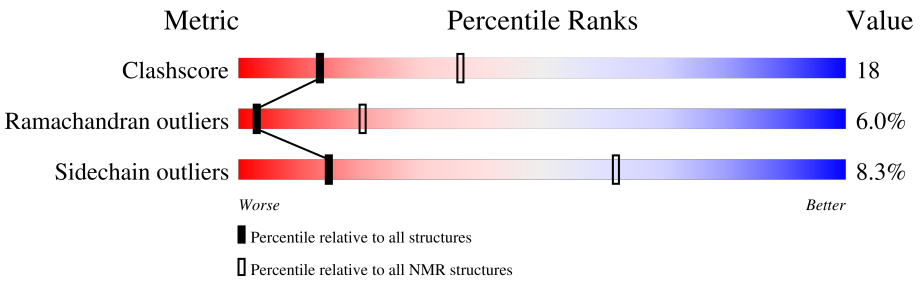
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.27
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.27

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	117	

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 5 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:12-A:103 (92)	0.23	5

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. No single-model clusters were found.

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

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 1736 atoms, of which 869 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Tjp2 protein.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	117	1736	537	869	156	171	3	0

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	cloning artifact	UNP Q9Z0U1
A	2	SER	-	cloning artifact	UNP Q9Z0U1
A	3	SER	-	cloning artifact	UNP Q9Z0U1
A	4	GLY	-	cloning artifact	UNP Q9Z0U1
A	5	SER	-	cloning artifact	UNP Q9Z0U1
A	6	SER	-	cloning artifact	UNP Q9Z0U1
A	7	GLY	-	cloning artifact	UNP Q9Z0U1
A	112	SER	-	cloning artifact	UNP Q9Z0U1
A	113	GLY	-	cloning artifact	UNP Q9Z0U1
A	114	PRO	-	cloning artifact	UNP Q9Z0U1
A	115	SER	-	cloning artifact	UNP Q9Z0U1
A	116	SER	-	cloning artifact	UNP Q9Z0U1
A	117	GLY	-	cloning artifact	UNP Q9Z0U1

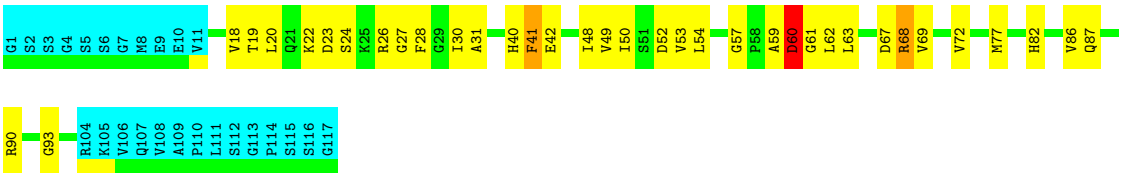
4 Residue-property plots (i)

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: Tjp2 protein

Chain A: 48% 28% • • 21%



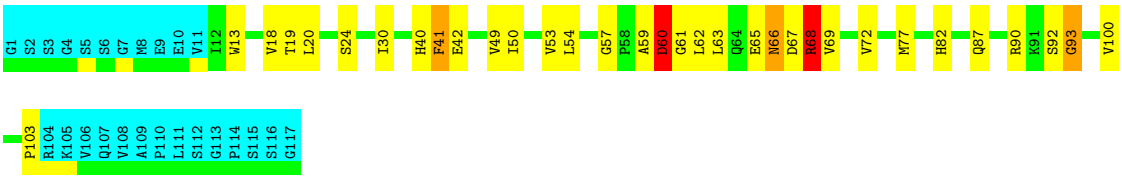
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

• Molecule 1: Tjp2 protein

Chain A: 50% 24% • • 21%



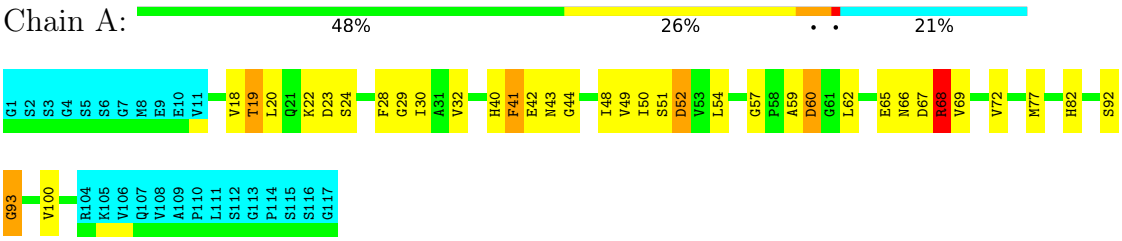
4.2.2 Score per residue for model 2

• Molecule 1: Tjp2 protein

Chain A: 51% 24% • • 21%

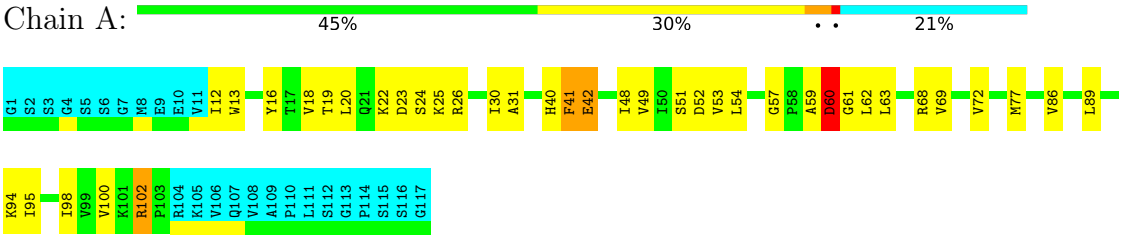
4.2.3 Score per residue for model 3

- Molecule 1: Tjp2 protein



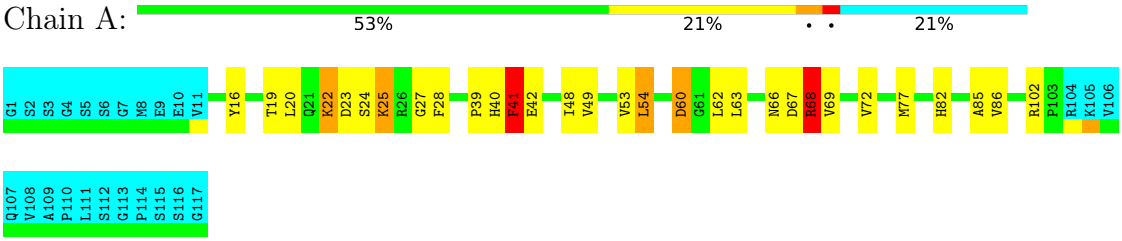
4.2.4 Score per residue for model 4

- Molecule 1: Tjp2 protein



4.2.5 Score per residue for model 5 (medoid)

- Molecule 1: Tjp2 protein



4.2.6 Score per residue for model 6

- Molecule 1: Tjp2 protein

Chain A: 50% 24% • • 21%

G1 S2 S3 G4 S5 S6 G7 M8 E9 E10 V11 T12 W13 T19 L20 Q21 K22 D23 S24 I30 A31 R36 H40 F41 E42 V49 I50 S51 D52 V53 L54 G57 P58 A59 D60 G61 L62 L63 R68 V72 M77 V80 L81 H82 S83 F84 A85 V86 Q87

G93 P103 R104 K105 V106 Q107 V108 A109 P110 L111 S112 G113 P114 S115 S116 G117

4.2.7 Score per residue for model 7

- Molecule 1: Tjp2 protein

Chain A: 50% 24% • • 21%

G1 S2 S3 G4 S5 S6 G7 M8 E9 E10 V11 Y16 T17 V18 T19 L20 Q21 K22 D23 S24 K25 R26 G27 F28 G29 I30 A31 H40 F41 E42 N43 G44 E45 V49 D52 V53 L54 A59 D60 G61 L62 L63 Q64 E65 N66 D67 R68 V69 M77 Q87 R90

R104 K105 V106 Q107 V108 A109 P110 L111 S112 G113 P114 S115 S116 G117

4.2.8 Score per residue for model 8

- Molecule 1: Tjp2 protein

Chain A: 48% 26% • • 21%

G1 S2 S3 G4 S5 S6 G7 M8 E9 E10 V11 T12 W13 V18 T19 L20 Q21 K22 R26 G27 F28 G29 I30 A31 H40 F41 E42 E45 T46 S47 I48 V49 I50 S51 D52 V53 L54 G57 P58 A59 D60 G61 L62 N66 D67 R68 V69 V86 Q87 Q88 L89 R90

V100 R104 K105 V106 Q107 V108 A109 P110 L111 S112 G113 P114 S115 S116 G117

4.2.9 Score per residue for model 9

- Molecule 1: Tjp2 protein

Chain A: 49% 24% • • 21%

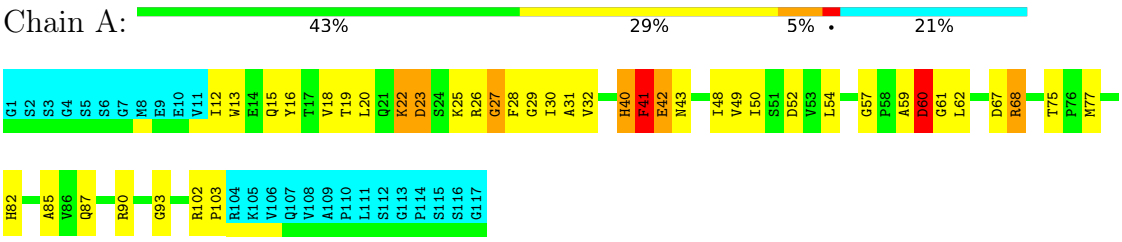
G1 S2 S3 G4 S5 S6 G7 M8 E9 E10 V11 Y16 T17 V18 T19 L20 Q21 K22 D23 R26 G27 F28 G29 I30 A31 H40 F41 E42 N43 G44 I48 V49 I50 S51 D52 A59 D60 G61 L62 L63 D67 R68 V69 V72 M77 H82 V86 Q87 R90

R102 P103 R104 K105 V106 Q107 V108 A109 P110 L111 S112 G113 P114 S115 S116 G117

4.2.10 Score per residue for model 10

- Molecule 1: Tjp2 protein

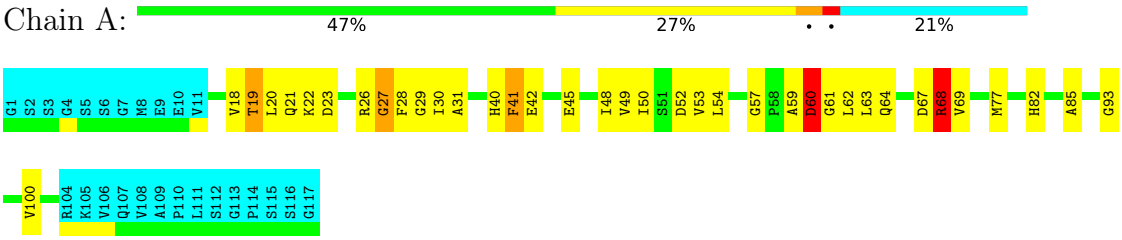
Chain A:



4.2.11 Score per residue for model 11

- Molecule 1: Tjp2 protein

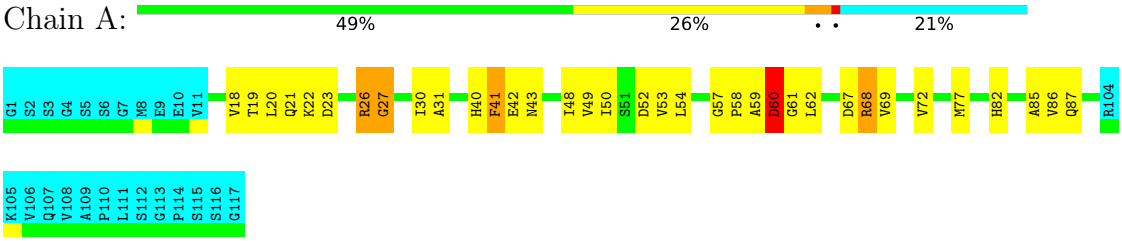
Chain A:



4.2.12 Score per residue for model 12

- Molecule 1: Tjp2 protein

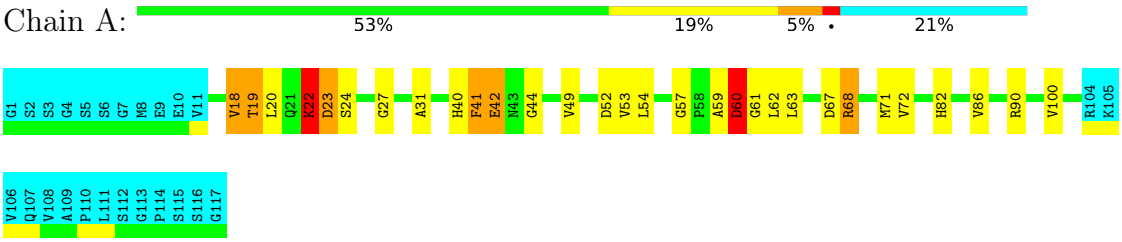
Chain A:



4.2.13 Score per residue for model 13

- Molecule 1: Tjp2 protein

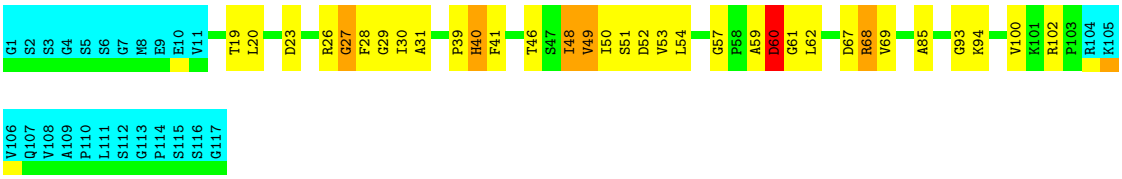
Chain A:



4.2.14 Score per residue for model 14

- Molecule 1: Tjp2 protein

Chain A: 50% 23% • • 21%



4.2.15 Score per residue for model 15

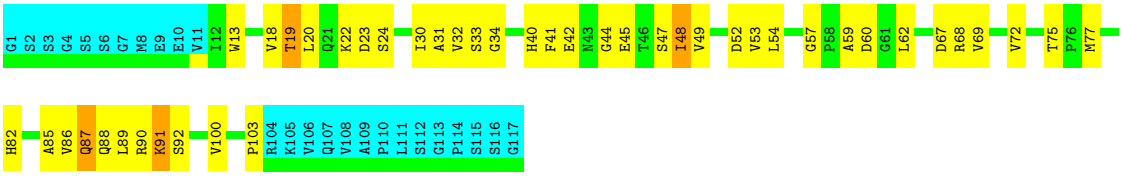
- Molecule 1: Tjp2 protein

Chain A: 45% 27% • • 21%

4.2.16 Score per residue for model 16

- Molecule 1: Tjp2 protein

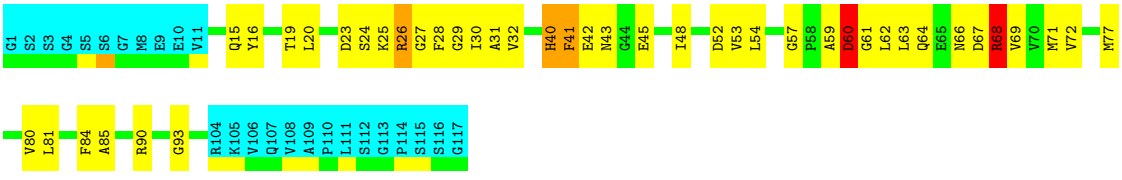
Chain A: 41% 34% • 21%



4.2.17 Score per residue for model 17

- Molecule 1: Tjp2 protein

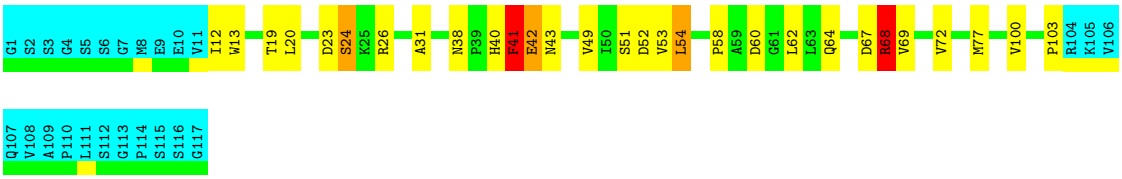
Chain A: 42% 32% • • 21%



4.2.18 Score per residue for model 18

- Molecule 1: Tjp2 protein

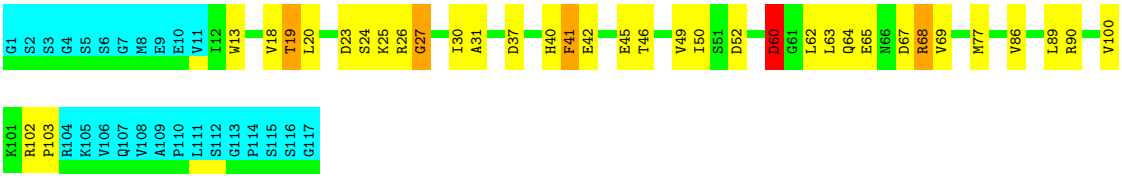
Chain A: 54% 21% • • 21%



4.2.19 Score per residue for model 19

- Molecule 1: Tjp2 protein

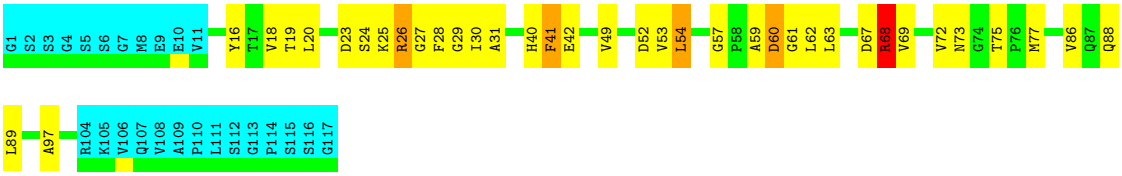
Chain A: 49% 26% • • 21%



4.2.20 Score per residue for model 20

- Molecule 1: Tjp2 protein

Chain A: 47% 27% • • 21%



5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *torsion angle dynamics*.

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *structures with the least restraint violation, target function*.

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

Software name	Classification	Version
CYANA	structure solution	2.0.17
CYANA	refinement	2.0.17

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	702	707	705	25±4
All	All	14040	14140	14100	509

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 18.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:67:ASP:OD2	1:A:100:VAL:HG21	0.87	1.69	13	3
1:A:18:VAL:HG21	1:A:62:LEU:O	0.83	1.73	12	9
1:A:54:LEU:HD23	1:A:57:GLY:HA3	0.80	1.51	15	12
1:A:20:LEU:HD23	1:A:62:LEU:CD1	0.77	2.10	7	19
1:A:41:PHE:CE2	1:A:49:VAL:HG21	0.76	2.16	14	1
1:A:82:HIS:O	1:A:86:VAL:HG12	0.75	1.82	12	5
1:A:20:LEU:HD23	1:A:62:LEU:HD12	0.75	1.58	7	19
1:A:31:ALA:HB3	1:A:52:ASP:HB2	0.68	1.66	2	8
1:A:54:LEU:HD23	1:A:57:GLY:CA	0.67	2.19	16	4
1:A:48:ILE:HG12	1:A:85:ALA:HB2	0.67	1.64	14	1
1:A:31:ALA:HB3	1:A:52:ASP:CB	0.67	2.20	20	11
1:A:63:LEU:HD22	1:A:100:VAL:HG21	0.66	1.66	4	1
1:A:26:ARG:O	1:A:27:GLY:C	0.66	2.33	10	8
1:A:20:LEU:HD11	1:A:63:LEU:HD12	0.65	1.68	6	9

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:31:ALA:HB3	1:A:52:ASP:HB3	0.64	1.70	12	9
1:A:30:ILE:HB	1:A:50:ILE:HD12	0.63	1.70	10	9
1:A:19:THR:O	1:A:62:LEU:HD13	0.62	1.95	11	8
1:A:63:LEU:CD2	1:A:100:VAL:HG21	0.61	2.24	4	1
1:A:48:ILE:O	1:A:69:VAL:HG23	0.61	1.96	14	5
1:A:58:PRO:O	1:A:62:LEU:HD12	0.61	1.95	12	2
1:A:30:ILE:CG2	1:A:59:ALA:CB	0.61	2.79	16	1
1:A:67:ASP:OD2	1:A:100:VAL:HG11	0.61	1.96	8	2
1:A:48:ILE:HG13	1:A:85:ALA:HB2	0.57	1.76	17	8
1:A:22:LYS:HE2	1:A:28:PHE:CZ	0.56	2.36	8	3
1:A:20:LEU:HD11	1:A:63:LEU:CD1	0.56	2.30	20	5
1:A:30:ILE:HG23	1:A:59:ALA:HB2	0.56	1.77	4	3
1:A:54:LEU:HD13	1:A:57:GLY:HA3	0.56	1.76	20	2
1:A:20:LEU:HD23	1:A:62:LEU:HD13	0.54	1.78	17	5
1:A:69:VAL:HG12	1:A:77:MET:HG3	0.54	1.80	20	8
1:A:48:ILE:CG1	1:A:85:ALA:HB2	0.54	2.33	14	1
1:A:63:LEU:CD2	1:A:100:VAL:CG2	0.54	2.86	4	1
1:A:72:VAL:HG23	1:A:77:MET:HG3	0.53	1.79	9	12
1:A:22:LYS:HE3	1:A:28:PHE:CZ	0.53	2.38	7	2
1:A:82:HIS:CD2	1:A:82:HIS:C	0.53	2.82	12	10
1:A:22:LYS:CE	1:A:28:PHE:CZ	0.52	2.93	9	4
1:A:28:PHE:O	1:A:30:ILE:HG23	0.52	2.05	14	3
1:A:32:VAL:HB	1:A:48:ILE:HG23	0.52	1.82	16	4
1:A:40:HIS:ND1	1:A:40:HIS:C	0.52	2.63	14	10
1:A:40:HIS:O	1:A:40:HIS:CG	0.51	2.63	17	1
1:A:13:TRP:CH2	1:A:68:ARG:NH2	0.51	2.79	18	1
1:A:40:HIS:O	1:A:41:PHE:C	0.51	2.48	7	19
1:A:41:PHE:O	1:A:42:GLU:C	0.51	2.49	11	17
1:A:87:GLN:O	1:A:90:ARG:CG	0.50	2.60	16	8
1:A:75:THR:HG21	1:A:88:GLN:NE2	0.50	2.21	16	1
1:A:66:ASN:O	1:A:68:ARG:CZ	0.50	2.60	7	2
1:A:26:ARG:HG2	1:A:54:LEU:HD22	0.50	1.82	12	1
1:A:72:VAL:HG22	1:A:98:ILE:HG23	0.50	1.84	4	1
1:A:26:ARG:O	1:A:29:GLY:N	0.50	2.44	10	1
1:A:48:ILE:CD1	1:A:48:ILE:N	0.49	2.75	14	1
1:A:59:ALA:O	1:A:61:GLY:N	0.49	2.45	14	15
1:A:41:PHE:O	1:A:44:GLY:N	0.49	2.46	3	3
1:A:65:GLU:O	1:A:66:ASN:ND2	0.49	2.45	3	2
1:A:67:ASP:CG	1:A:100:VAL:HG21	0.49	2.28	13	1
1:A:34:GLY:O	1:A:48:ILE:HD13	0.49	2.07	16	1
1:A:67:ASP:O	1:A:69:VAL:N	0.48	2.46	18	7

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:30:ILE:CG2	1:A:59:ALA:HB2	0.48	2.38	4	1
1:A:41:PHE:O	1:A:43:ASN:N	0.48	2.46	10	3
1:A:48:ILE:O	1:A:48:ILE:HG22	0.48	2.07	14	1
1:A:13:TRP:CZ3	1:A:68:ARG:HG3	0.48	2.44	16	1
1:A:53:VAL:O	1:A:54:LEU:C	0.48	2.52	15	16
1:A:67:ASP:O	1:A:68:ARG:C	0.48	2.52	18	10
1:A:80:VAL:HB	1:A:84:PHE:CD1	0.48	2.44	6	1
1:A:60:ASP:CG	1:A:60:ASP:O	0.48	2.53	18	7
1:A:32:VAL:HG23	1:A:82:HIS:CE1	0.47	2.44	10	2
1:A:42:GLU:CD	1:A:43:ASN:N	0.47	2.68	3	2
1:A:40:HIS:O	1:A:40:HIS:ND1	0.47	2.48	7	1
1:A:48:ILE:CG1	1:A:85:ALA:CB	0.47	2.92	14	1
1:A:45:GLU:OE1	1:A:68:ARG:NE	0.47	2.47	16	1
1:A:28:PHE:O	1:A:29:GLY:C	0.47	2.53	14	9
1:A:51:SER:C	1:A:52:ASP:OD1	0.47	2.54	3	4
1:A:86:VAL:O	1:A:89:LEU:N	0.47	2.48	19	6
1:A:51:SER:O	1:A:52:ASP:CG	0.47	2.54	18	5
1:A:13:TRP:CZ3	1:A:103:PRO:HD3	0.46	2.45	10	8
1:A:16:TYR:CE1	1:A:102:ARG:HB2	0.46	2.45	4	3
1:A:67:ASP:OD2	1:A:100:VAL:CG2	0.46	2.60	16	1
1:A:64:GLN:CG	1:A:67:ASP:OD1	0.46	2.63	7	2
1:A:25:LYS:O	1:A:26:ARG:C	0.46	2.53	20	3
1:A:25:LYS:O	1:A:27:GLY:N	0.46	2.49	17	2
1:A:50:ILE:CG1	1:A:67:ASP:OD1	0.46	2.64	8	3
1:A:41:PHE:CD2	1:A:45:GLU:HB3	0.46	2.46	7	2
1:A:48:ILE:HG21	1:A:85:ALA:HB1	0.45	1.88	14	1
1:A:60:ASP:O	1:A:60:ASP:CG	0.45	2.55	10	10
1:A:66:ASN:O	1:A:68:ARG:NH1	0.45	2.49	5	4
1:A:53:VAL:HG13	1:A:59:ALA:HB1	0.45	1.88	4	1
1:A:22:LYS:CE	1:A:28:PHE:CE2	0.45	2.99	5	1
1:A:47:SER:O	1:A:48:ILE:C	0.45	2.55	16	1
1:A:45:GLU:O	1:A:46:THR:C	0.45	2.54	19	1
1:A:39:PRO:O	1:A:40:HIS:C	0.45	2.55	2	3
1:A:48:ILE:O	1:A:69:VAL:CG2	0.45	2.63	14	1
1:A:26:ARG:O	1:A:27:GLY:O	0.45	2.34	15	5
1:A:80:VAL:HB	1:A:84:PHE:CD2	0.45	2.46	17	1
1:A:16:TYR:CD1	1:A:16:TYR:N	0.45	2.85	9	5
1:A:73:ASN:OD1	1:A:97:ALA:N	0.45	2.49	20	1
1:A:75:THR:HG21	1:A:88:GLN:CD	0.45	2.32	20	1
1:A:20:LEU:HD13	1:A:28:PHE:HB3	0.45	1.89	9	1
1:A:63:LEU:HD22	1:A:67:ASP:OD2	0.45	2.11	19	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:67:ASP:O	1:A:67:ASP:CG	0.44	2.54	3	3
1:A:82:HIS:O	1:A:82:HIS:CD2	0.44	2.70	3	1
1:A:63:LEU:HD22	1:A:100:VAL:CG2	0.44	2.40	4	1
1:A:41:PHE:C	1:A:43:ASN:N	0.44	2.71	17	1
1:A:23:ASP:N	1:A:27:GLY:HA2	0.44	2.27	15	7
1:A:23:ASP:O	1:A:24:SER:C	0.44	2.56	19	2
1:A:40:HIS:O	1:A:42:GLU:N	0.43	2.51	7	3
1:A:30:ILE:HG23	1:A:59:ALA:CB	0.43	2.42	16	1
1:A:26:ARG:O	1:A:28:PHE:N	0.43	2.51	7	1
1:A:47:SER:CB	1:A:68:ARG:HD2	0.43	2.43	16	1
1:A:26:ARG:HG3	1:A:54:LEU:HD11	0.43	1.91	18	1
1:A:86:VAL:O	1:A:87:GLN:C	0.43	2.57	16	2
1:A:71:MET:HG2	1:A:72:VAL:N	0.43	2.28	17	2
1:A:75:THR:HG22	1:A:77:MET:CE	0.43	2.43	10	2
1:A:94:LYS:O	1:A:95:ILE:HD13	0.43	2.14	4	1
1:A:15:GLN:C	1:A:16:TYR:CD1	0.43	2.92	10	2
1:A:100:VAL:HG21	1:A:102:ARG:NH2	0.43	2.29	14	1
1:A:64:GLN:O	1:A:67:ASP:CG	0.42	2.58	11	2
1:A:31:ALA:CB	1:A:52:ASP:HB3	0.42	2.41	20	1
1:A:90:ARG:CZ	1:A:90:ARG:HB2	0.42	2.44	17	4
1:A:92:SER:O	1:A:93:GLY:C	0.42	2.58	3	2
1:A:90:ARG:HB2	1:A:90:ARG:CZ	0.42	2.45	13	1
1:A:21:GLN:O	1:A:27:GLY:O	0.42	2.38	11	2
1:A:67:ASP:O	1:A:67:ASP:OD1	0.42	2.38	14	1
1:A:25:LYS:C	1:A:27:GLY:N	0.41	2.74	5	2
1:A:60:ASP:OD1	1:A:60:ASP:C	0.41	2.58	10	1
1:A:82:HIS:CD2	1:A:82:HIS:O	0.41	2.73	5	1
1:A:39:PRO:HB3	1:A:46:THR:HG22	0.41	1.91	14	1
1:A:13:TRP:CE2	1:A:103:PRO:HB3	0.41	2.51	15	2
1:A:64:GLN:HG3	1:A:65:GLU:N	0.41	2.30	19	1
1:A:67:ASP:HB3	1:A:100:VAL:HG21	0.41	1.92	11	1
1:A:87:GLN:O	1:A:91:LYS:HG3	0.41	2.15	16	1
1:A:54:LEU:HD22	1:A:57:GLY:N	0.40	2.31	20	1
1:A:67:ASP:HB2	1:A:100:VAL:HG21	0.40	1.94	1	1
1:A:91:LYS:O	1:A:92:SER:C	0.40	2.59	16	1
1:A:73:ASN:OD1	1:A:97:ALA:O	0.40	2.38	20	1
1:A:22:LYS:O	1:A:23:ASP:O	0.40	2.38	6	2
1:A:34:GLY:O	1:A:48:ILE:CD1	0.40	2.69	16	1
1:A:40:HIS:C	1:A:40:HIS:ND1	0.40	2.75	5	1
1:A:81:LEU:O	1:A:84:PHE:N	0.40	2.55	17	1
1:A:38:ASN:O	1:A:38:ASN:CG	0.40	2.59	18	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:67:ASP:HB3	1:A:100:VAL:HG11	0.40	1.92	18	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	92/117 (79%)	72±2 (79±2%)	14±3 (15±3%)	6±1 (6±1%)	3	20
All	All	1840/2340 (79%)	1447 (79%)	283 (15%)	110 (6%)	3	20

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	60	ASP	18
1	A	68	ARG	18
1	A	41	PHE	16
1	A	23	ASP	13
1	A	24	SER	11
1	A	93	GLY	9
1	A	27	GLY	9
1	A	22	LYS	8
1	A	42	GLU	3
1	A	44	GLY	2
1	A	26	ARG	2
1	A	48	ILE	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	77/96 (80%)	71±1 (92±2%)	6±1 (8±2%)	15 62
All	All	1540/1920 (80%)	1412 (92%)	128 (8%)	15 62

All 26 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	19	THR	20
1	A	49	VAL	19
1	A	60	ASP	18
1	A	68	ARG	13
1	A	41	PHE	7
1	A	18	VAL	6
1	A	40	HIS	6
1	A	54	LEU	6
1	A	22	LYS	5
1	A	102	ARG	4
1	A	25	LYS	4
1	A	87	GLN	4
1	A	42	GLU	2
1	A	45	GLU	2
1	A	66	ASN	1
1	A	52	ASP	1
1	A	24	SER	1
1	A	36	ARG	1
1	A	43	ASN	1
1	A	47	SER	1
1	A	26	ARG	1
1	A	48	ILE	1
1	A	94	LYS	1
1	A	33	SER	1
1	A	91	LYS	1
1	A	37	ASP	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