Full wwPDB NMR Structure Validation Report (i)

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PDB ID	:	2B9Z
Title	:	Solution structure of FHV B2, a viral suppressor of RNAi
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Deposited on	:	2005-10-13

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:

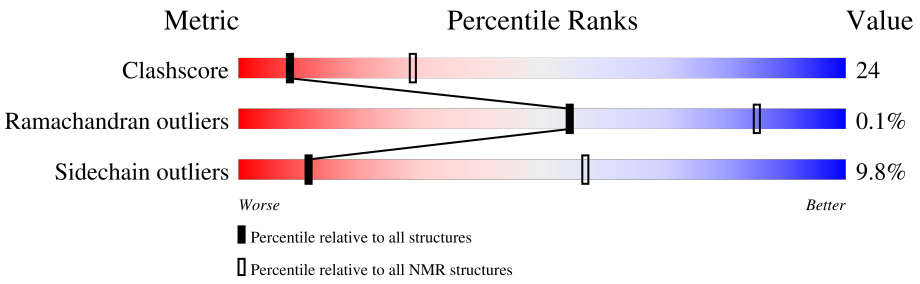
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.27
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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	74	49% 42% • 7%
1	B	74	54% 39% • •

2 Ensemble composition and analysis (i)

This entry contains 10 models. Model 7 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	**Residue range (total)**	**Backbone RMSD (Å)**	**Medoid model**
1	A:3-A:71, B:2-B:72 (140)	0.28	7

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

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

The models can be grouped into 2 clusters. No single-model clusters were found.

Cluster number	Models
1	1, 2, 6, 7, 8, 9, 10
2	3, 4, 5

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2294 atoms, of which 1172 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called B2 protein.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	74	1147	349	586	100	107	5	0
1	B	74	1147	349	586	100	107	5	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	cloning artifact	UNP P68831
A	0	ALA	-	cloning artifact	UNP P68831
A	53	GLY	SER	SEE REMARK 999	UNP P68831
B	-1	GLY	-	cloning artifact	UNP P68831
B	0	ALA	-	cloning artifact	UNP P68831
B	53	GLY	SER	SEE REMARK 999	UNP P68831

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: B2 protein

Chain A: 49% 42% • 7%

• Molecule 1: B2 protein

Chain B: 54% 39% • •

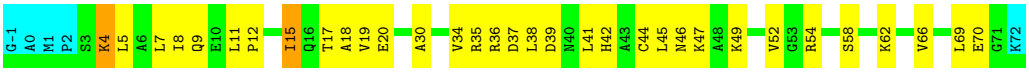
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

• Molecule 1: B2 protein

Chain A: 49% 42% • 7%



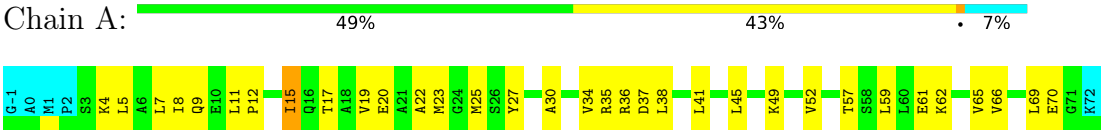
• Molecule 1: B2 protein

Chain B: 47% 45% • •

4.2.2 Score per residue for model 2

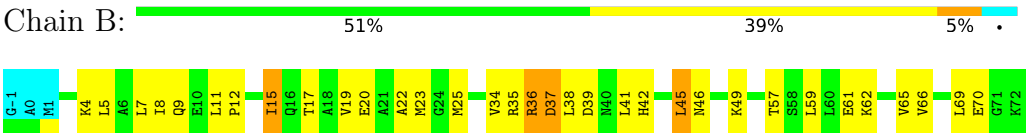
• Molecule 1: B2 protein

Chain A:



• Molecule 1: B2 protein

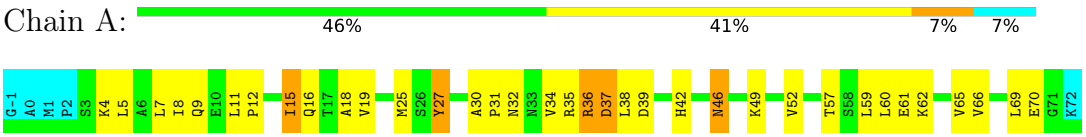
Chain B:



4.2.3 Score per residue for model 3

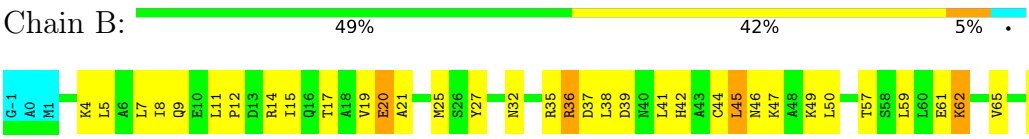
• Molecule 1: B2 protein

Chain A:



• Molecule 1: B2 protein

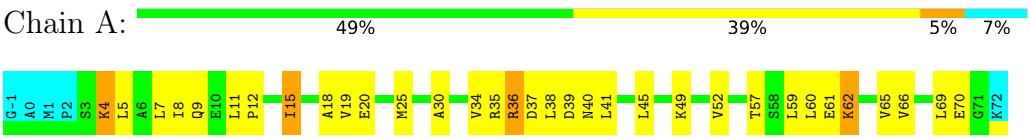
Chain B:



4.2.4 Score per residue for model 4

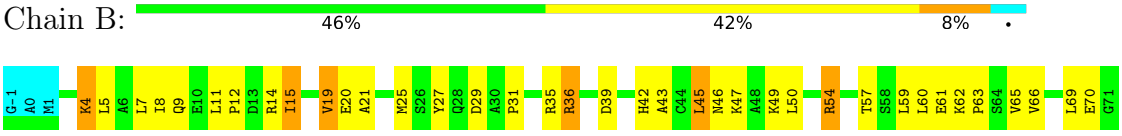
• Molecule 1: B2 protein

Chain A:



• Molecule 1: B2 protein

Chain B:



K72

4.2.5 Score per residue for model 5

- Molecule 1: B2 protein

Chain A: 55% 32% 5% 7%

G-1 A0 M1 P2 S3 K4 L5 A6 L7 I8 Q9 E10 L11 P12 I15 Q16 T17 A18 V19 E20 A21 A22 M25 S26 Y27 V34 R35 R36 D37 L38 D39 H42 N46 K49 M55 V56 T57 L60 E61 L69 E70 G71 K72

- Molecule 1: B2 protein

Chain B: 51% 39% 5% •

G-1 A0 M1 K4 L5 A6 L7 I8 Q9 E10 L11 P12 I15 V19 E20 A21 M25 S26 Y27 P31 N32 N33 V34 R35 R36 D37 L38 D39 H42 N46 K49 L50 M55 S58 L59 L60 E61 K62 V65 V66 L69 E70 G71 K72

4.2.6 Score per residue for model 6

- Molecule 1: B2 protein

Chain A: 50% 35% 8% 7%

G-1 A0 M1 P2 S3 K4 L5 A6 L7 I8 Q9 E10 L11 P12 D13 R14 I15 A18 V19 M23 G24 M25 N32 N33 V34 R35 R36 D37 L38 D39 N40 L41 H42 L45 K49 L50 S58 L59 L60 E61 K62 V66 L69 E70 G71 K72

- Molecule 1: B2 protein

Chain B: 54% 35% 7% •

G-1 A0 M1 K4 L5 A6 L7 I8 L11 P12 D13 R14 I15 E20 M25 S26 Y27 N32 R35 R36 D39 N40 L41 H42 A43 C44 L45 N46 K47 A48 K49 L50 T51 V52 S58 E61 K62 V65 V66 L69 E70 G71 K72

4.2.7 Score per residue for model 7 (medoid)

- Molecule 1: B2 protein

Chain A: 50% 36% 7% 7%

G-1 A0 M1 P2 S3 K4 L5 A6 L7 I8 Q9 E10 L11 P12 I15 Q16 T17 A18 V19 E20 M25 S26 Y27 N32 R35 R36 D37 L38 D39 N40 L41 H42 A43 C44 L45 N46 K47 A48 K49 V52 T57 S58 L59 L60 E61 V65 L69 K72

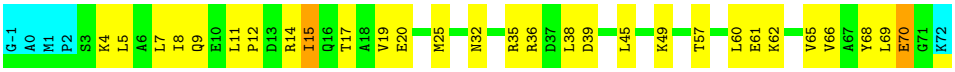
- Molecule 1: B2 protein

Chain B: 47% 39% 9% •

4.2.8 Score per residue for model 8

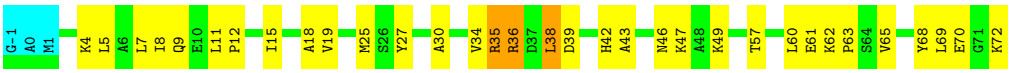
- Molecule 1: B2 protein

Chain A:



- Molecule 1: B2 protein

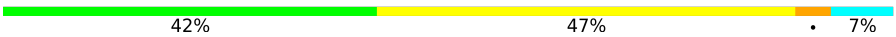
Chain B:



4.2.9 Score per residue for model 9

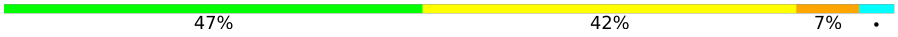
- Molecule 1: B2 protein

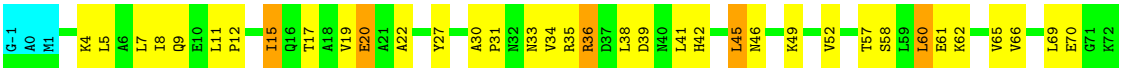
Chain A:



- Molecule 1: B2 protein

Chain B:

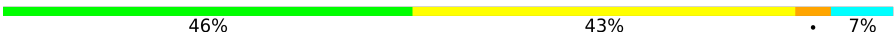


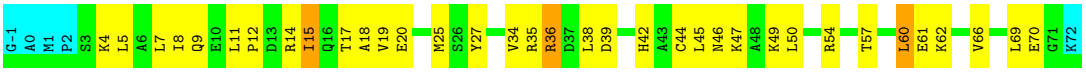


4.2.10 Score per residue for model 10

- Molecule 1: B2 protein

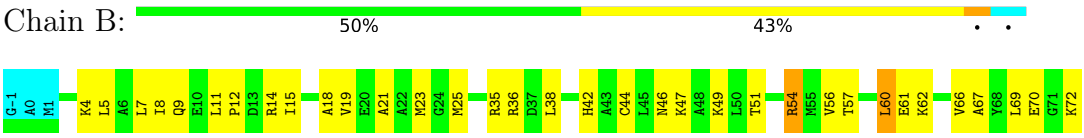
Chain A:





• Molecule 1: B2 protein

Chain B:



5 Refinement protocol and experimental data overview (i)

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

Of the 100 calculated structures, 10 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	1.2
CNS	structure solution	1.1
CNS	refinement	1.1

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	527	547	546	30±4
1	B	544	567	566	29±5
All	All	10710	11140	11120	526

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 24.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:19:VAL:HG21	1:A:49:LYS:HD3	0.86	1.48	10	3
1:B:19:VAL:HG21	1:B:49:LYS:HD3	0.81	1.53	10	3
1:B:4:LYS:HG3	1:B:69:LEU:HA	0.80	1.52	9	2
1:A:25:MET:SD	1:B:4:LYS:HB2	0.75	2.20	4	1
1:A:4:LYS:HG2	1:A:69:LEU:HA	0.75	1.55	4	1
1:B:34:VAL:O	1:B:38:LEU:HD23	0.73	1.83	8	2
1:A:25:MET:SD	1:B:4:LYS:HB3	0.73	2.24	10	4
1:B:5:LEU:HA	1:B:8:ILE:HG13	0.73	1.61	9	10
1:A:52:VAL:HG13	1:B:45:LEU:HD21	0.72	1.59	1	4
1:A:59:LEU:HD12	1:A:65:VAL:HG11	0.71	1.63	7	3
1:A:19:VAL:HG12	1:A:23:MET:SD	0.70	2.26	2	1
1:A:4:LYS:HB3	1:B:25:MET:SD	0.69	2.27	2	3
1:A:5:LEU:HA	1:A:8:ILE:HG13	0.69	1.63	9	10

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:14:ARG:HB3	1:B:18:ALA:HA	0.69	1.63	10	2
1:A:4:LYS:HB2	1:A:69:LEU:HA	0.68	1.65	10	3
1:B:5:LEU:HA	1:B:8:ILE:CG1	0.68	2.19	9	10
1:B:43:ALA:O	1:B:47:LYS:HG3	0.67	1.89	6	1
1:A:44:CYS:HA	1:A:47:LYS:HE2	0.67	1.65	9	2
1:A:42:HIS:O	1:A:46:ASN:HB3	0.67	1.89	10	6
1:B:59:LEU:HD12	1:B:65:VAL:HG11	0.67	1.65	3	4
1:A:19:VAL:HG12	1:A:23:MET:HG2	0.66	1.66	6	2
1:A:25:MET:SD	1:B:4:LYS:HE3	0.66	2.31	10	1
1:A:15:ILE:HG23	1:A:49:LYS:HG2	0.66	1.66	2	10
1:B:19:VAL:CG2	1:B:49:LYS:HD3	0.65	2.21	8	4
1:B:5:LEU:HG	1:B:8:ILE:HD11	0.65	1.68	3	5
1:A:5:LEU:HA	1:A:8:ILE:CG1	0.65	2.20	9	10
1:B:42:HIS:O	1:B:46:ASN:HB3	0.65	1.91	4	10
1:A:4:LYS:CG	1:A:69:LEU:HA	0.65	2.22	4	1
1:A:58:SER:O	1:A:62:LYS:HE2	0.64	1.93	1	3
1:A:68:TYR:HE2	1:B:27:TYR:HA	0.64	1.52	8	1
1:B:67:ALA:HB1	1:B:72:LYS:HB2	0.63	1.67	10	1
1:A:5:LEU:HG	1:A:8:ILE:HD11	0.62	1.69	10	8
1:B:4:LYS:HG2	1:B:69:LEU:O	0.62	1.93	4	1
1:A:68:TYR:CE2	1:B:27:TYR:HA	0.61	2.29	8	1
1:A:22:ALA:O	1:A:25:MET:HB2	0.61	1.96	2	1
1:A:34:VAL:HG13	1:B:65:VAL:HA	0.61	1.70	9	1
1:A:58:SER:O	1:A:61:GLU:HG3	0.61	1.94	6	1
1:B:41:LEU:O	1:B:45:LEU:HB2	0.61	1.95	3	6
1:A:41:LEU:O	1:A:45:LEU:HB2	0.61	1.96	7	6
1:B:36:ARG:O	1:B:39:ASP:HB3	0.60	1.97	7	5
1:A:8:ILE:HG12	1:A:69:LEU:HD13	0.59	1.75	1	6
1:B:19:VAL:HG11	1:B:45:LEU:HD12	0.59	1.73	3	1
1:A:23:MET:HE1	1:A:42:HIS:HA	0.59	1.74	6	1
1:A:44:CYS:O	1:A:47:LYS:HG2	0.59	1.97	10	1
1:B:4:LYS:HG3	1:B:69:LEU:O	0.59	1.98	1	1
1:A:19:VAL:CG2	1:A:49:LYS:HD3	0.58	2.27	10	3
1:B:8:ILE:HG12	1:B:69:LEU:HD13	0.58	1.74	4	8
1:B:44:CYS:HA	1:B:47:LYS:HE2	0.58	1.75	1	2
1:A:57:THR:O	1:A:61:GLU:HG2	0.58	1.98	4	8
1:B:4:LYS:CG	1:B:69:LEU:HA	0.58	2.28	9	2
1:B:46:ASN:O	1:B:50:LEU:HG	0.58	1.98	5	1
1:A:7:LEU:O	1:A:11:LEU:HG	0.58	1.98	7	10
1:B:19:VAL:HG12	1:B:49:LYS:HD3	0.57	1.75	1	3
1:A:8:ILE:HG12	1:A:69:LEU:HD22	0.57	1.76	4	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:17:THR:HA	1:A:20:GLU:CG	0.57	2.30	9	2
1:A:25:MET:SD	1:B:4:LYS:HE2	0.57	2.39	8	1
1:A:35:ARG:O	1:A:38:LEU:HB2	0.56	2.00	10	1
1:A:68:TYR:CZ	1:B:30:ALA:HB2	0.56	2.35	8	1
1:B:36:ARG:HD2	1:B:39:ASP:OD2	0.56	1.99	9	2
1:A:45:LEU:HD11	1:B:52:VAL:HG13	0.56	1.77	1	3
1:A:35:ARG:HA	1:A:38:LEU:HD23	0.56	1.77	2	4
1:B:35:ARG:HA	1:B:38:LEU:HD23	0.56	1.78	1	4
1:B:57:THR:O	1:B:61:GLU:HG2	0.56	2.01	9	6
1:A:19:VAL:HG21	1:A:45:LEU:HB3	0.56	1.78	4	1
1:A:36:ARG:O	1:A:39:ASP:HB3	0.55	2.01	3	6
1:B:4:LYS:HG3	1:B:69:LEU:CA	0.55	2.31	9	1
1:B:4:LYS:HB2	1:B:69:LEU:HA	0.55	1.78	10	2
1:A:32:ASN:HA	1:A:35:ARG:HG2	0.55	1.79	7	2
1:A:4:LYS:HG3	1:A:69:LEU:O	0.55	2.02	1	1
1:A:4:LYS:CD	1:A:69:LEU:HA	0.55	2.31	4	1
1:A:4:LYS:HG2	1:A:69:LEU:CA	0.54	2.29	4	1
1:B:27:TYR:CD1	1:B:35:ARG:HB2	0.53	2.37	6	2
1:A:27:TYR:CD2	1:A:35:ARG:HB2	0.53	2.39	5	3
1:B:36:ARG:HA	1:B:39:ASP:HB2	0.53	1.81	9	4
1:A:27:TYR:HD2	1:A:35:ARG:HB2	0.53	1.63	7	4
1:A:27:TYR:H	1:B:4:LYS:HE3	0.53	1.63	3	1
1:A:4:LYS:HB2	1:A:69:LEU:O	0.53	2.03	5	1
1:A:38:LEU:HD11	1:B:69:LEU:HD21	0.53	1.81	3	3
1:A:41:LEU:HD21	1:B:59:LEU:HB2	0.53	1.80	4	1
1:A:27:TYR:HB2	1:A:38:LEU:HD23	0.53	1.79	5	1
1:A:34:VAL:O	1:A:37:ASP:HB2	0.53	2.04	1	3
1:A:69:LEU:HD21	1:B:38:LEU:HD11	0.53	1.81	1	2
1:A:4:LYS:HE2	1:B:25:MET:SD	0.53	2.44	5	2
1:B:46:ASN:HA	1:B:49:LYS:NZ	0.53	2.17	5	1
1:B:58:SER:O	1:B:62:LYS:HE2	0.53	2.03	1	2
1:A:44:CYS:HA	1:A:47:LYS:CD	0.53	2.33	7	2
1:B:27:TYR:CD2	1:B:35:ARG:HB2	0.52	2.39	9	2
1:A:68:TYR:CE2	1:B:30:ALA:HB2	0.52	2.40	8	1
1:B:22:ALA:O	1:B:25:MET:HB2	0.52	2.05	2	1
1:B:35:ARG:O	1:B:38:LEU:HB2	0.52	2.04	8	1
1:A:17:THR:HA	1:A:20:GLU:HG2	0.51	1.83	2	5
1:B:43:ALA:O	1:B:47:LYS:HB2	0.51	2.05	8	2
1:A:52:VAL:HG13	1:B:45:LEU:HD11	0.51	1.81	7	2
1:A:36:ARG:HA	1:A:39:ASP:HB2	0.51	1.80	9	2
1:A:68:TYR:CE2	1:B:31:PRO:HD3	0.51	2.41	9	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:19:VAL:CG1	1:A:49:LYS:HD3	0.51	2.36	7	1
1:B:27:TYR:HD2	1:B:35:ARG:HB2	0.51	1.64	4	3
1:B:27:TYR:HD1	1:B:35:ARG:HB2	0.51	1.65	6	2
1:A:36:ARG:HD2	1:A:39:ASP:OD2	0.51	2.06	10	2
1:A:7:LEU:HG	1:B:21:ALA:HB1	0.50	1.83	1	5
1:A:27:TYR:CD1	1:A:35:ARG:HB2	0.50	2.41	3	1
1:B:32:ASN:O	1:B:35:ARG:HG2	0.50	2.07	3	2
1:A:9:GLN:O	1:A:12:PRO:HD2	0.50	2.06	8	10
1:A:35:ARG:HA	1:A:38:LEU:HB2	0.50	1.82	9	1
1:A:66:VAL:O	1:A:70:GLU:HG2	0.50	2.06	9	6
1:B:27:TYR:HB2	1:B:38:LEU:HG	0.50	1.83	9	2
1:A:25:MET:SD	1:A:27:TYR:HB3	0.50	2.46	5	2
1:B:65:VAL:O	1:B:69:LEU:HG	0.50	2.06	6	2
1:A:65:VAL:O	1:A:69:LEU:HG	0.50	2.06	8	1
1:A:5:LEU:HB2	1:A:69:LEU:CB	0.50	2.36	4	10
1:B:7:LEU:O	1:B:11:LEU:HG	0.49	2.06	8	10
1:A:31:PRO:O	1:A:35:ARG:HG2	0.49	2.07	3	1
1:A:44:CYS:HA	1:A:47:LYS:HD3	0.49	1.84	7	1
1:B:9:GLN:O	1:B:12:PRO:HD2	0.49	2.05	7	9
1:A:27:TYR:HB2	1:A:38:LEU:HG	0.49	1.83	9	1
1:B:66:VAL:O	1:B:70:GLU:HG2	0.49	2.07	6	8
1:A:34:VAL:O	1:A:38:LEU:HD23	0.49	2.07	10	2
1:B:19:VAL:CG1	1:B:49:LYS:HD3	0.49	2.38	4	3
1:B:44:CYS:HA	1:B:47:LYS:CD	0.49	2.38	7	1
1:B:5:LEU:HB2	1:B:69:LEU:HB3	0.49	1.83	9	1
1:B:30:ALA:O	1:B:35:ARG:HD3	0.48	2.09	9	1
1:A:4:LYS:HG3	1:B:25:MET:SD	0.48	2.48	4	1
1:A:5:LEU:HB2	1:A:69:LEU:HB3	0.48	1.86	4	1
1:B:27:TYR:O	1:B:35:ARG:HB2	0.48	2.08	8	1
1:A:33:ASN:HA	1:A:36:ARG:HB2	0.48	1.84	9	1
1:A:37:ASP:OD2	1:B:62:LYS:HE3	0.47	2.09	3	1
1:A:4:LYS:HE2	1:B:25:MET:CG	0.47	2.39	3	1
1:A:18:ALA:HA	1:B:14:ARG:HB3	0.47	1.86	10	5
1:A:25:MET:CG	1:B:4:LYS:HE2	0.47	2.38	5	1
1:A:27:TYR:HD1	1:A:35:ARG:HB2	0.47	1.67	3	1
1:A:62:LYS:O	1:A:66:VAL:HG23	0.47	2.10	9	4
1:B:58:SER:O	1:B:62:LYS:HE3	0.47	2.10	7	1
1:B:19:VAL:HB	1:B:49:LYS:HD3	0.47	1.85	7	1
1:A:36:ARG:O	1:A:39:ASP:HB2	0.47	2.10	9	1
1:A:15:ILE:O	1:A:19:VAL:HG23	0.46	2.09	2	1
1:B:15:ILE:HG23	1:B:49:LYS:HG2	0.46	1.87	2	5

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:50:LEU:O	1:B:54:ARG:HD2	0.46	2.10	4	1
1:A:22:ALA:O	1:A:25:MET:HB3	0.46	2.10	5	1
1:B:47:LYS:O	1:B:51:THR:HG23	0.46	2.10	10	2
1:A:19:VAL:HG12	1:A:45:LEU:HG	0.46	1.86	8	2
1:A:36:ARG:HA	1:A:39:ASP:CB	0.46	2.39	7	3
1:B:5:LEU:HB2	1:B:69:LEU:CB	0.46	2.39	9	8
1:A:50:LEU:O	1:A:54:ARG:HB3	0.46	2.10	10	1
1:B:36:ARG:HA	1:B:39:ASP:CB	0.46	2.41	1	1
1:B:27:TYR:HD2	1:B:35:ARG:CB	0.46	2.23	9	1
1:A:30:ALA:O	1:A:35:ARG:HD3	0.46	2.09	1	1
1:B:62:LYS:O	1:B:66:VAL:HG23	0.46	2.10	2	4
1:A:46:ASN:HA	1:A:49:LYS:NZ	0.46	2.26	5	1
1:A:52:VAL:CG1	1:B:45:LEU:HD11	0.46	2.41	3	1
1:B:44:CYS:O	1:B:47:LYS:HG2	0.46	2.11	7	2
1:A:32:ASN:O	1:A:35:ARG:HG2	0.46	2.10	6	1
1:B:54:ARG:HA	1:B:57:THR:OG1	0.46	2.11	1	1
1:A:59:LEU:CD1	1:A:69:LEU:HD11	0.46	2.41	4	1
1:A:32:ASN:HA	1:A:35:ARG:NE	0.45	2.25	9	2
1:A:11:LEU:HB2	1:A:12:PRO:HD3	0.45	1.89	2	8
1:B:17:THR:HA	1:B:20:GLU:CG	0.45	2.41	3	1
1:B:34:VAL:O	1:B:38:LEU:HD22	0.45	2.11	5	1
1:B:15:ILE:HD12	1:B:49:LYS:HG2	0.45	1.87	7	1
1:A:30:ALA:HB1	1:A:34:VAL:HG23	0.45	1.88	3	3
1:A:15:ILE:HD13	1:A:19:VAL:HG23	0.45	1.88	6	2
1:A:27:TYR:O	1:A:35:ARG:HB2	0.45	2.12	10	1
1:A:66:VAL:HA	1:A:69:LEU:HD12	0.45	1.88	4	2
1:A:62:LYS:HG3	1:A:65:VAL:HB	0.44	1.87	4	1
1:B:32:ASN:O	1:B:36:ARG:HG2	0.44	2.13	1	1
1:A:34:VAL:O	1:A:38:LEU:HD22	0.44	2.12	5	2
1:A:19:VAL:HG12	1:A:49:LYS:HD3	0.44	1.89	7	1
1:B:58:SER:O	1:B:61:GLU:HG3	0.44	2.13	6	2
1:A:12:PRO:O	1:A:16:GLN:HG3	0.44	2.13	3	1
1:B:17:THR:HA	1:B:20:GLU:HG2	0.44	1.90	3	4
1:A:8:ILE:HD13	1:A:60:LEU:HD21	0.44	1.89	9	1
1:B:11:LEU:HB2	1:B:12:PRO:HD3	0.43	1.89	9	7
1:A:32:ASN:O	1:A:36:ARG:HD2	0.43	2.14	6	1
1:B:34:VAL:O	1:B:37:ASP:HB2	0.43	2.12	1	2
1:A:17:THR:O	1:A:20:GLU:HG3	0.43	2.13	1	1
1:B:4:LYS:HB2	1:B:69:LEU:O	0.43	2.14	5	1
1:B:72:LYS:HD2	1:B:72:LYS:C	0.43	2.34	7	1
1:B:46:ASN:HA	1:B:49:LYS:CE	0.43	2.44	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:44:CYS:HA	1:A:47:LYS:HG2	0.42	1.91	7	1
1:B:33:ASN:HA	1:B:36:ARG:HB2	0.42	1.90	9	1
1:B:32:ASN:HA	1:B:35:ARG:NE	0.42	2.29	5	1
1:A:32:ASN:O	1:A:36:ARG:HD3	0.42	2.14	8	1
1:A:11:LEU:HD11	1:B:22:ALA:HB2	0.42	1.91	9	1
1:A:68:TYR:CZ	1:B:31:PRO:HD3	0.42	2.49	9	1
1:B:56:VAL:O	1:B:60:LEU:HB2	0.42	2.14	10	1
1:A:27:TYR:N	1:B:4:LYS:HE3	0.42	2.27	3	1
1:B:27:TYR:CB	1:B:38:LEU:HG	0.42	2.44	7	1
1:B:5:LEU:CA	1:B:8:ILE:HG13	0.42	2.40	9	1
1:B:19:VAL:HG21	1:B:45:LEU:HD12	0.42	1.90	9	1
1:B:54:ARG:O	1:B:57:THR:HB	0.42	2.14	10	1
1:A:30:ALA:HB1	1:A:34:VAL:CG2	0.42	2.45	2	1
1:B:59:LEU:CD2	1:B:65:VAL:HG11	0.42	2.44	4	1
1:B:35:ARG:HA	1:B:38:LEU:CD2	0.42	2.44	3	1
1:B:62:LYS:HG3	1:B:65:VAL:HB	0.41	1.92	4	1
1:B:60:LEU:HA	1:B:66:VAL:HG22	0.41	1.92	9	1
1:A:60:LEU:HA	1:A:66:VAL:HG22	0.41	1.91	10	1
1:B:31:PRO:O	1:B:35:ARG:HG2	0.41	2.15	4	2
1:B:19:VAL:HG21	1:B:49:LYS:CE	0.41	2.45	5	1
1:A:4:LYS:HB3	1:B:25:MET:HE2	0.41	1.92	8	1
1:A:19:VAL:HG21	1:A:49:LYS:CD	0.41	2.33	10	1
1:B:59:LEU:HD11	1:B:69:LEU:HD11	0.41	1.93	4	1
1:A:4:LYS:HD2	1:B:25:MET:SD	0.41	2.55	2	1
1:A:32:ASN:CA	1:A:35:ARG:HG2	0.41	2.45	7	1
1:B:44:CYS:HA	1:B:47:LYS:HG2	0.41	1.91	7	1
1:A:15:ILE:CG2	1:A:49:LYS:HG2	0.41	2.44	8	1
1:B:11:LEU:N	1:B:12:PRO:HD2	0.41	2.30	6	1
1:A:50:LEU:O	1:A:50:LEU:HD12	0.40	2.16	6	1
1:A:19:VAL:HB	1:A:49:LYS:HD3	0.40	1.94	1	1
1:B:62:LYS:HA	1:B:63:PRO:HD3	0.40	1.79	4	3
1:B:16:GLN:O	1:B:20:GLU:HB3	0.40	2.15	7	1
1:B:70:GLU:HB2	1:B:72:LYS:HG2	0.40	1.93	8	1
1:A:19:VAL:CG1	1:A:45:LEU:HG	0.40	2.46	8	1
1:A:65:VAL:HA	1:B:34:VAL:HG13	0.40	1.93	9	1
1:A:4:LYS:HE3	1:B:25:MET:SD	0.40	2.56	1	1
1:B:19:VAL:O	1:B:23:MET:HG2	0.40	2.17	2	1
1:A:4:LYS:HE3	1:A:68:TYR:CE2	0.40	2.52	8	1
1:A:15:ILE:O	1:A:19:VAL:HG13	0.40	2.17	10	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	69/74 (93%)	68±1 (98±2%)	1±1 (2±2%)	0±0 (0±0%)	54	85
1	B	70/74 (95%)	69±1 (99±1%)	1±1 (1±1%)	0±0 (0±0%)	54	85
All	All	1390/1480 (94%)	1367 (98%)	21 (2%)	2 (0%)	54	85

All 2 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	27	TYR	1
1	B	27	TYR	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	58/61 (95%)	53±2 (91±4%)	5±2 (9±4%)	13	60
1	B	60/61 (98%)	54±2 (89±3%)	6±2 (11±3%)	10	54
All	All	1180/1220 (97%)	1064 (90%)	116 (10%)	11	57

All 40 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	15	ILE	10
1	B	15	ILE	10
1	B	36	ARG	9
1	A	36	ARG	7
1	A	60	LEU	7

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Mol	Chain	Res	Type	Models (Total)
1	B	20	GLU	6
1	B	45	LEU	6
1	B	60	LEU	5
1	A	62	LYS	4
1	B	37	ASP	4
1	B	62	LYS	4
1	A	4	LYS	3
1	A	37	ASP	3
1	A	46	ASN	3
1	B	4	LYS	2
1	B	19	VAL	2
1	B	50	LEU	2
1	A	20	GLU	2
1	B	54	ARG	2
1	A	38	LEU	2
1	A	70	GLU	2
1	B	46	ASN	2
1	B	61	GLU	2
1	A	54	ARG	1
1	A	40	ASN	1
1	B	29	ASP	1
1	A	55	MET	1
1	B	55	MET	1
1	A	13	ASP	1
1	A	50	LEU	1
1	A	61	GLU	1
1	A	19	VAL	1
1	A	45	LEU	1
1	B	70	GLU	1
1	B	72	LYS	1
1	B	35	ARG	1
1	B	38	LEU	1
1	B	68	TYR	1
1	A	42	HIS	1
1	B	23	MET	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