



Full wwPDB NMR Structure Validation Report ⓘ

Feb 12, 2017 – 10:03 pm GMT

PDB ID : 2HFI
Title : Solution NMR Structure of Protein yppE from *Bacillus subtilis*. Northeast Structural Genomics Consortium Target SR213
Authors : Liu, G.; Singarapu, K.K.; Parish, D.; Eletsky, A.; Xu, D.; Sukumaran, D.; Ho, C.K.; Fang, Y.; Cunningham, K.; Ma, L.-C.; Xiao, R.; Liu, J.; Baran, M.; Swapna, G.V.T.; Acton, T.B.; Rost, B.; Montelione, G.T.; Szyperski, T.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2006-06-23

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

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

with specific help available everywhere you see the ⓘ symbol.

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

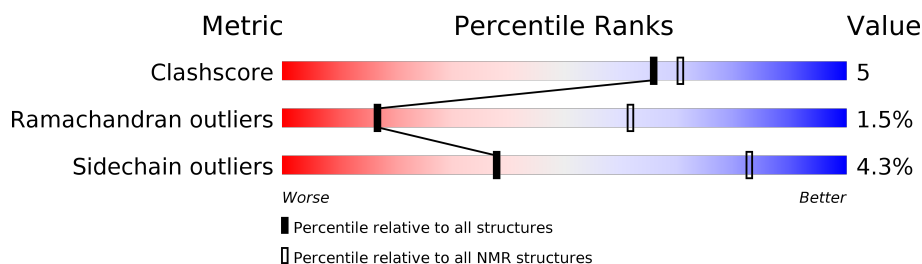
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 78%.

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 ≥ 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	131	

2 Ensemble composition and analysis

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

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:6-A:25, A:41-A:120 (100)	0.50	20

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2031 atoms, of which 1013 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Hypothetical protein yppE.

Mol	Chain	Residues	Atoms						Trace
1	A	123	Total	C	H	N	O	S	0
			2031	642	1013	177	196	3	

There are 8 discrepancies between the modelled and reference sequences:

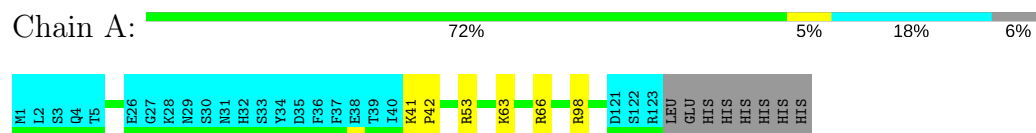
Chain	Residue	Modelled	Actual	Comment	Reference
A	124	LEU	-	CLONING ARTIFACT	UNP P50833
A	125	GLU	-	CLONING ARTIFACT	UNP P50833
A	126	HIS	-	EXPRESSION TAG	UNP P50833
A	127	HIS	-	EXPRESSION TAG	UNP P50833
A	128	HIS	-	EXPRESSION TAG	UNP P50833
A	129	HIS	-	EXPRESSION TAG	UNP P50833
A	130	HIS	-	EXPRESSION TAG	UNP P50833
A	131	HIS	-	EXPRESSION TAG	UNP P50833

4 Residue-property plots

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA and DNA chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: Hypothetical protein yppE

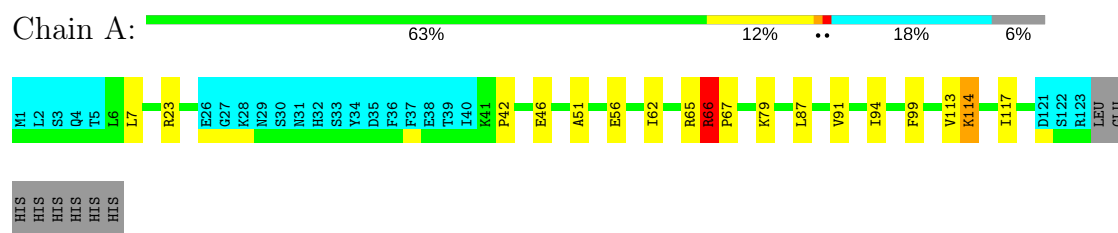


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

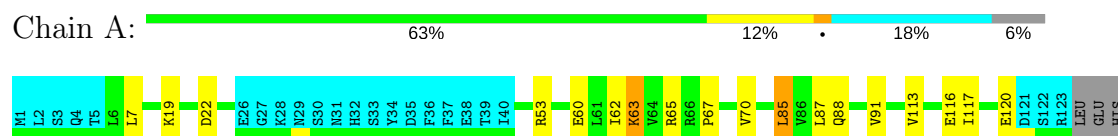
4.2.1 Score per residue for model 1

- Molecule 1: Hypothetical protein yppE



4.2.2 Score per residue for model 2

- Molecule 1: Hypothetical protein yppE



HIS
HIS
HIS
HIS

4.2.3 Score per residue for model 3

- Molecule 1: Hypothetical protein yppE

Chain A: 66% 8% 18% 6%

M1 L2 L3 S3 Q4 T5 E26 G27 K28 N29 S30 N31 H32 S33 Y34 D35 F36 F37 E38 T39 I40 K41 P42 R53 K63 V64 R65 R66 Q74 V78 E84 L87 Q88 V91 E120 D121 S122 R123 LEU GLU HIS HIS HIS HIS HIS HIS

4.2.4 Score per residue for model 4

- Molecule 1: Hypothetical protein yppE

Chain A: 63% 11% 18% 6%

M1 L2 L3 S3 Q4 T5 E8 E13 M13 I14 E15 K19 E26 G27 K28 N29 S30 N31 H32 S33 Y34 D35 F36 F37 E38 T39 I40 K41 P42 E46 E60 K63 V64 R65 R66 E84 Q88 H95 K96 K97 R98 L110 D121 S122 R123 LEU GLU HIS HIS HIS HIS HIS

HIS
HIS
HIS

4.2.5 Score per residue for model 5

- Molecule 1: Hypothetical protein yppE

Chain A: 65% 11% 18% 6%

M1 L2 L3 S3 Q4 T5 L6 L7 E8 M13 R23 E26 G27 K28 N29 S30 N31 H32 S33 Y34 D35 F36 F37 E38 T39 I40 K41 P67 V70 Q74 V78 I94 H95 K96 K97 R98 L110 K114 D121 S122 R123 LEU GLU HIS HIS HIS HIS HIS HIS

4.2.6 Score per residue for model 6

- Molecule 1: Hypothetical protein yppE

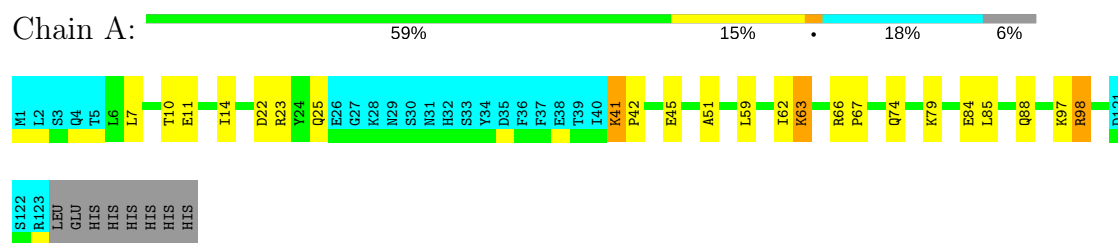
Chain A: 56% 21% 18% 6%

M1 L2 L3 S3 Q4 T5 L6 L7 T10 E11 Q12 M13 I14 D22 R23 K28 N29 S30 N31 H32 S33 Y34 D35 F36 F37 E38 T39 I40 A51 A52 R53 L59 E60 K63 V64 R65 R66 P67 K72 K79 E84 Q88 H93 K96 K100

E104 L110 D121 S122 R123 LEU GLU HIS HIS HIS HIS HIS HIS

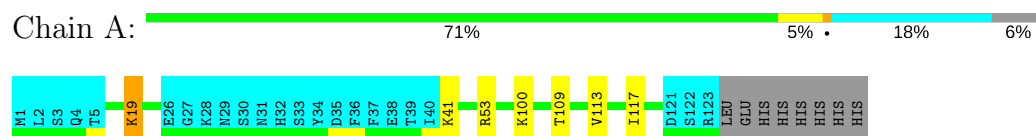
4.2.7 Score per residue for model 7

- Molecule 1: Hypothetical protein yppE



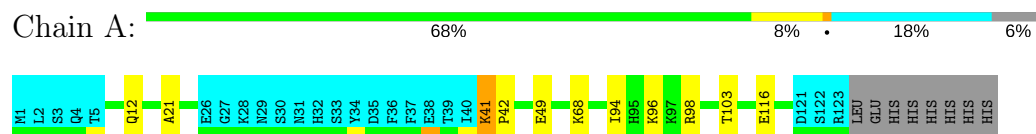
4.2.8 Score per residue for model 8

- Molecule 1: Hypothetical protein yppE



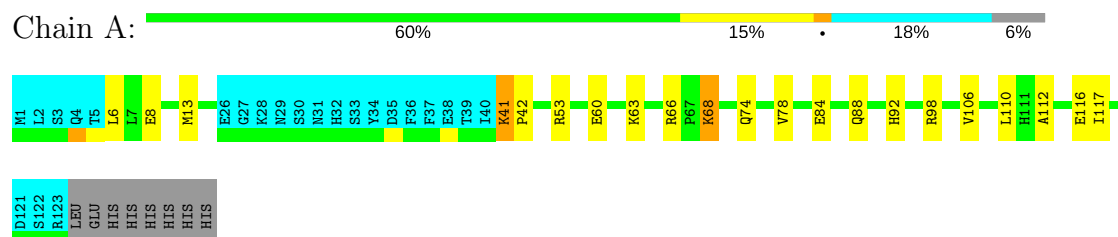
4.2.9 Score per residue for model 9

- Molecule 1: Hypothetical protein yppE



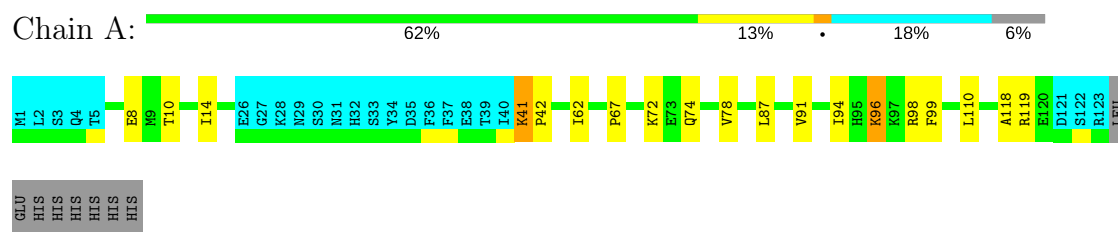
4.2.10 Score per residue for model 10

- Molecule 1: Hypothetical protein yppE



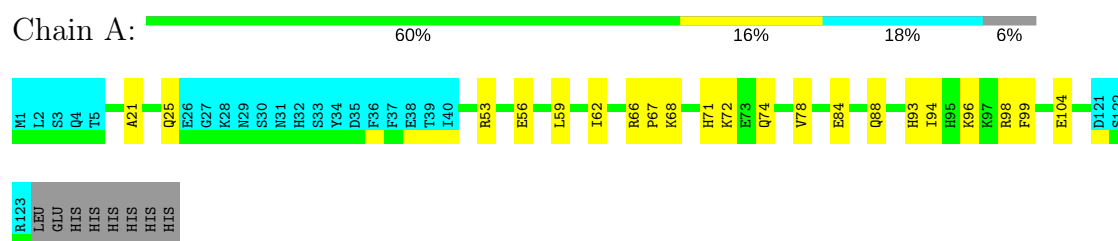
4.2.11 Score per residue for model 11

- Molecule 1: Hypothetical protein yppE



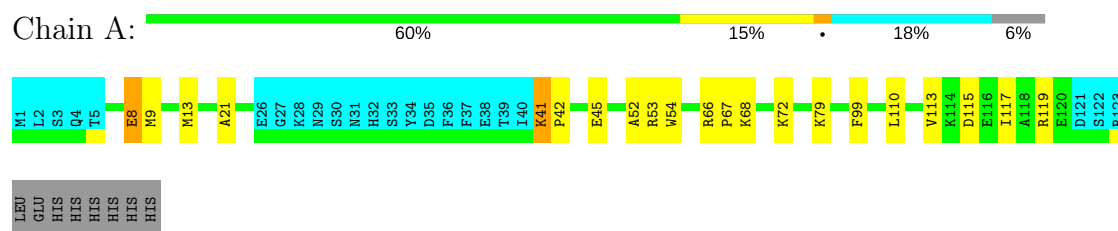
4.2.12 Score per residue for model 12

- Molecule 1: Hypothetical protein yppE



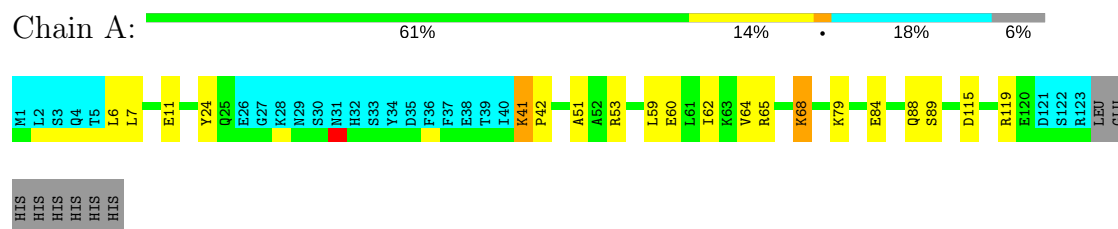
4.2.13 Score per residue for model 13

- Molecule 1: Hypothetical protein yppE



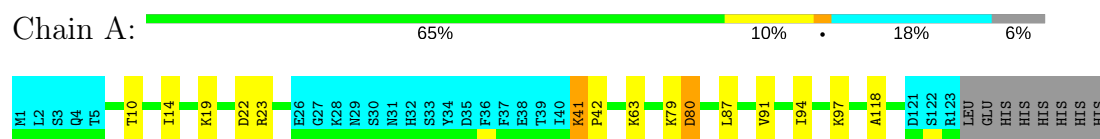
4.2.14 Score per residue for model 14

- Molecule 1: Hypothetical protein yppE



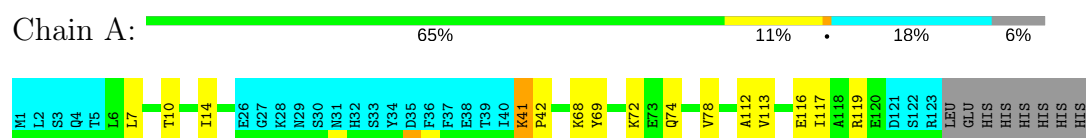
4.2.15 Score per residue for model 15

- Molecule 1: Hypothetical protein yppE



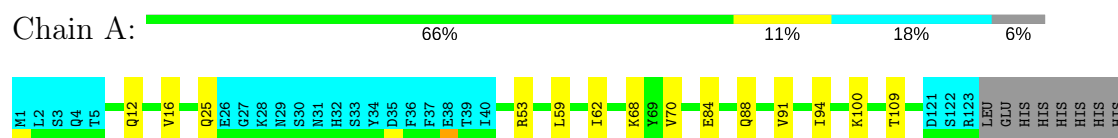
4.2.16 Score per residue for model 16

- Molecule 1: Hypothetical protein yppE



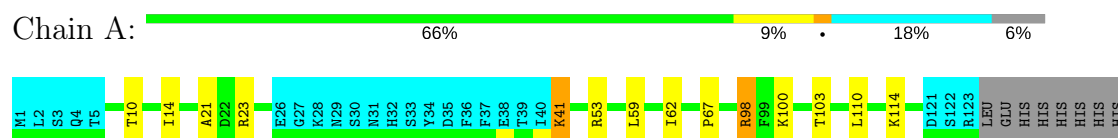
4.2.17 Score per residue for model 17

- Molecule 1: Hypothetical protein yppE



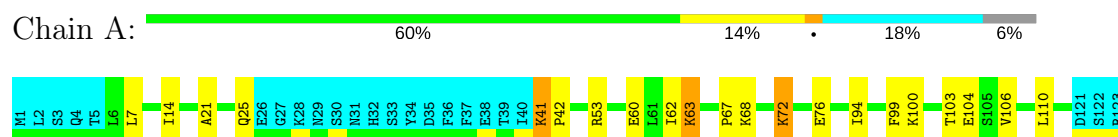
4.2.18 Score per residue for model 18

- Molecule 1: Hypothetical protein yppE



4.2.19 Score per residue for model 19

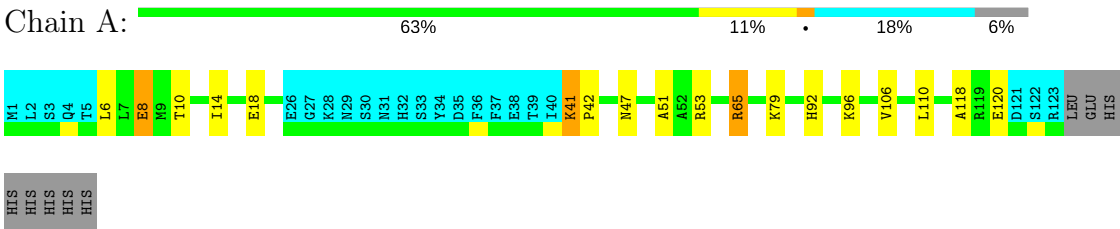
- Molecule 1: Hypothetical protein yppE



LEU
GLU
HIS
HIS
HIS
HIS
HIS
HIS

4.2.20 Score per residue for model 20 (medoid)

- Molecule 1: Hypothetical protein yppE



5 Refinement protocol and experimental data overview

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

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

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

Software name	Classification	Version
CYANA	structure solution	2.1
CNS	refinement	1.1

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	BMRB entry 7227
Number of chemical shift lists	1
Total number of shifts	1302
Number of shifts mapped to atoms	1302
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	78%

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	829	842	837	8±3
All	All	16580	16840	16740	158

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 5.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:41:LYS:HE2	1:A:41:LYS:HA	0.76	1.58	7	1
1:A:65:ARG:HD2	1:A:120:GLU:HA	0.72	1.59	3	1
1:A:14:ILE:HG12	1:A:110:LEU:HB3	0.71	1.61	11	3
1:A:94:ILE:HD11	1:A:98:ARG:HB2	0.65	1.66	11	1
1:A:22:ASP:O	1:A:25:GLN:HG3	0.64	1.93	6	1
1:A:94:ILE:HG22	1:A:98:ARG:HD3	0.63	1.69	12	1
1:A:85:LEU:HA	1:A:88:GLN:HE21	0.62	1.54	7	1
1:A:62:ILE:HB	1:A:67:PRO:HD3	0.61	1.72	1	2
1:A:72:LYS:O	1:A:76:GLU:HG2	0.61	1.95	19	1
1:A:51:ALA:HB1	1:A:79:LYS:HG3	0.61	1.73	20	3
1:A:94:ILE:HG12	1:A:99:PHE:HB2	0.60	1.74	1	3
1:A:59:LEU:HD23	1:A:62:ILE:HD11	0.59	1.74	12	4
1:A:66:ARG:HG2	1:A:67:PRO:HD2	0.58	1.75	13	1
1:A:24:TYR:HE2	1:A:96:LYS:HA	0.58	1.58	6	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:52:ALA:HA	1:A:79:LYS:HE3	0.58	1.74	13	1
1:A:116:GLU:O	1:A:120:GLU:HG2	0.56	2.00	2	1
1:A:59:LEU:O	1:A:63:LYS:HG3	0.56	2.00	7	1
1:A:62:ILE:HG22	1:A:67:PRO:HG3	0.56	1.77	11	1
1:A:24:TYR:CD1	1:A:89:SER:HA	0.55	2.36	14	1
1:A:65:ARG:HD2	1:A:120:GLU:CA	0.55	2.31	3	1
1:A:62:ILE:HB	1:A:67:PRO:HG3	0.55	1.78	12	2
1:A:7:LEU:HG	1:A:11:GLU:HG2	0.55	1.76	6	1
1:A:60:GLU:O	1:A:63:LYS:HG2	0.54	2.03	10	1
1:A:115:ASP:O	1:A:119:ARG:HB2	0.54	2.03	13	2
1:A:68:LYS:HE3	1:A:116:GLU:OE1	0.53	2.04	9	1
1:A:66:ARG:HA	1:A:66:ARG:HE	0.53	1.63	1	1
1:A:60:GLU:O	1:A:63:LYS:HG3	0.52	2.04	2	3
1:A:68:LYS:HB3	1:A:116:GLU:HG2	0.52	1.80	16	1
1:A:84:GLU:O	1:A:88:GLN:HG2	0.52	2.04	14	4
1:A:22:ASP:O	1:A:25:GLN:HG2	0.52	2.04	7	1
1:A:65:ARG:O	1:A:66:ARG:HB2	0.51	2.05	4	2
1:A:24:TYR:HD1	1:A:89:SER:HA	0.51	1.65	14	1
1:A:24:TYR:CE2	1:A:96:LYS:HA	0.51	2.38	6	1
1:A:74:GLN:O	1:A:78:VAL:HG23	0.50	2.07	5	6
1:A:68:LYS:NZ	1:A:68:LYS:HB3	0.50	2.21	14	1
1:A:13:MET:SD	1:A:110:LEU:HD21	0.50	2.45	6	5
1:A:42:PRO:O	1:A:46:GLU:HG2	0.50	2.06	1	1
1:A:87:LEU:O	1:A:91:VAL:HG22	0.50	2.07	11	5
1:A:94:ILE:HB	1:A:98:ARG:HB2	0.50	1.82	5	1
1:A:106:VAL:O	1:A:110:LEU:HG	0.49	2.08	20	3
1:A:94:ILE:HD12	1:A:98:ARG:HB2	0.49	1.82	9	1
1:A:22:ASP:HA	1:A:25:GLN:CG	0.49	2.38	6	1
1:A:41:LYS:O	1:A:45:GLU:HG2	0.49	2.08	13	1
1:A:7:LEU:O	1:A:11:GLU:HG2	0.49	2.08	14	1
1:A:84:GLU:O	1:A:88:GLN:HG3	0.49	2.07	12	3
1:A:9:MET:SD	1:A:53:ARG:HD2	0.48	2.49	13	1
1:A:7:LEU:HD21	1:A:114:LYS:HG2	0.47	1.85	1	1
1:A:66:ARG:HG2	1:A:67:PRO:CD	0.47	2.39	13	1
1:A:98:ARG:HA	1:A:98:ARG:NE	0.47	2.23	18	1
1:A:41:LYS:HB3	1:A:42:PRO:HD3	0.47	1.87	15	8
1:A:113:VAL:O	1:A:117:ILE:HG13	0.47	2.09	2	5
1:A:68:LYS:HB2	1:A:68:LYS:NZ	0.47	2.24	10	2
1:A:19:LYS:N	1:A:19:LYS:HE3	0.47	2.25	8	1
1:A:97:LYS:HG3	1:A:98:ARG:N	0.46	2.24	7	1
1:A:51:ALA:HB1	1:A:79:LYS:HG2	0.46	1.87	6	2

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:62:ILE:HG22	1:A:67:PRO:HG2	0.46	1.87	18	1
1:A:10:THR:O	1:A:14:ILE:HG13	0.45	2.10	15	6
1:A:100:LYS:O	1:A:104:GLU:HG2	0.45	2.11	6	2
1:A:84:GLU:HG3	1:A:88:GLN:NE2	0.45	2.26	17	1
1:A:65:ARG:HA	1:A:65:ARG:NE	0.45	2.26	3	1
1:A:67:PRO:HB2	1:A:70:VAL:O	0.45	2.11	2	2
1:A:79:LYS:HD3	1:A:80:ASP:N	0.45	2.27	15	1
1:A:21:ALA:HB2	1:A:103:THR:HG21	0.44	1.89	18	3
1:A:85:LEU:HA	1:A:88:GLN:HG2	0.44	1.90	2	1
1:A:21:ALA:HA	1:A:99:PHE:CZ	0.44	2.47	12	1
1:A:68:LYS:HD3	1:A:68:LYS:N	0.44	2.27	10	1
1:A:69:TYR:HB2	1:A:112:ALA:HB1	0.44	1.89	16	1
1:A:62:ILE:HA	1:A:65:ARG:O	0.44	2.13	2	1
1:A:41:LYS:HB2	1:A:42:PRO:HD3	0.44	1.90	11	2
1:A:62:ILE:HD12	1:A:72:LYS:HE2	0.43	1.89	11	1
1:A:70:VAL:HG21	1:A:109:THR:HG23	0.43	1.89	17	1
1:A:19:LYS:HA	1:A:22:ASP:HB3	0.43	1.90	15	2
1:A:112:ALA:O	1:A:116:GLU:HG2	0.43	2.13	10	1
1:A:88:GLN:HA	1:A:91:VAL:HG22	0.43	1.91	17	1
1:A:10:THR:O	1:A:14:ILE:HG12	0.42	2.14	7	1
1:A:60:GLU:O	1:A:64:VAL:HG23	0.42	2.15	6	2
1:A:62:ILE:HG22	1:A:67:PRO:HB3	0.42	1.90	19	1
1:A:66:ARG:HG3	1:A:67:PRO:HD2	0.42	1.91	6	1
1:A:68:LYS:HD2	1:A:68:LYS:N	0.42	2.30	19	1
1:A:21:ALA:HA	1:A:99:PHE:CE2	0.42	2.50	13	1
1:A:9:MET:CE	1:A:54:TRP:HA	0.42	2.44	13	1
1:A:41:LYS:HB3	1:A:42:PRO:CD	0.42	2.45	10	2
1:A:53:ARG:HH21	1:A:56:GLU:CD	0.42	2.18	12	1
1:A:65:ARG:HD3	1:A:120:GLU:OE2	0.41	2.14	20	1
1:A:68:LYS:HD2	1:A:69:TYR:CE2	0.41	2.50	16	1
1:A:88:GLN:HG3	1:A:94:ILE:HD13	0.41	1.92	17	1
1:A:8:GLU:CD	1:A:8:GLU:H	0.41	2.18	10	1
1:A:96:LYS:HE3	1:A:96:LYS:HA	0.41	1.92	11	1
1:A:74:GLN:OE1	1:A:74:GLN:HA	0.41	2.15	7	1
1:A:66:ARG:HA	1:A:66:ARG:NE	0.41	2.30	1	1
1:A:8:GLU:H	1:A:8:GLU:CD	0.41	2.16	13	1
1:A:15:GLU:HG2	1:A:19:LYS:HD3	0.41	1.93	4	1
1:A:12:GLN:O	1:A:16:VAL:HG23	0.41	2.16	17	1
1:A:22:ASP:HA	1:A:25:GLN:HG3	0.41	1.93	6	1
1:A:59:LEU:HD21	1:A:72:LYS:HG3	0.41	1.93	6	1
1:A:88:GLN:OE1	1:A:93:HIS:HB3	0.40	2.17	6	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:109:THR:O	1:A:113:VAL:HG23	0.40	2.16	8	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	100/131 (76%)	95±2 (95±2%)	4±2 (4±2%)	2±1 (2±1%)	17	62
All	All	2000/2620 (76%)	1895 (95%)	75 (4%)	30 (2%)	17	62

All 7 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	41	LYS	15
1	A	96	LYS	5
1	A	6	LEU	4
1	A	66	ARG	2
1	A	92	HIS	2
1	A	68	LYS	1
1	A	95	HIS	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	88/118 (75%)	84±1 (96±1%)	4±1 (4±1%)	38	83
All	All	1760/2360 (75%)	1685 (96%)	75 (4%)	38	83

All 31 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	ARG	10
1	A	63	LYS	7
1	A	23	ARG	5
1	A	72	LYS	4
1	A	66	ARG	4
1	A	98	ARG	4
1	A	68	LYS	3
1	A	25	GLN	3
1	A	114	LYS	3
1	A	100	LYS	3
1	A	7	LEU	3
1	A	65	ARG	3
1	A	8	GLU	3
1	A	119	ARG	2
1	A	12	GLN	2
1	A	47	ASN	1
1	A	80	ASP	1
1	A	94	ILE	1
1	A	19	LYS	1
1	A	18	GLU	1
1	A	46	GLU	1
1	A	95	HIS	1
1	A	104	GLU	1
1	A	93	HIS	1
1	A	49	GLU	1
1	A	45	GLU	1
1	A	97	LYS	1
1	A	96	LYS	1
1	A	71	HIS	1
1	A	85	LEU	1
1	A	56	GLU	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

The completeness of assignment taking into account all chemical shift lists is 78% for the well-defined parts and 70% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 7227

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	1302
Number of shifts mapped to atoms	1302
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	109	-0.56 ± 0.14	Should be applied
$^{13}\text{C}_\beta$	106	0.32 ± 0.11	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	95	0.74 ± 0.31	Should be applied

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 78%, i.e. 1031 atoms were assigned a chemical shift out of a possible 1321. 19 out of 19 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	377/496 (76%)	190/198 (96%)	97/200 (48%)	90/98 (92%)
Sidechain	596/723 (82%)	370/421 (88%)	221/268 (82%)	5/34 (15%)

Continued on next page...

Continued from previous page...

	Total	¹ H	¹³ C	¹⁵ N
Aromatic	58/102 (57%)	36/52 (69%)	21/39 (54%)	1/11 (9%)
Overall	1031/1321 (78%)	596/671 (89%)	339/507 (67%)	96/143 (67%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 70%, i.e. 1127 atoms were assigned a chemical shift out of a possible 1605. 19 out of 20 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	407/611 (67%)	205/244 (84%)	107/246 (43%)	95/121 (79%)
Sidechain	642/858 (75%)	400/501 (80%)	237/316 (75%)	5/41 (12%)
Aromatic	78/136 (57%)	50/70 (71%)	27/53 (51%)	1/13 (8%)
Overall	1127/1605 (70%)	655/815 (80%)	371/615 (60%)	101/175 (58%)

7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition.

Random coil index (RCI) for chain A:

