



Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 02:55 PM BST

PDB ID : 1GM1
Title : SECOND PDZ DOMAIN (PDZ2) OF PTP-BL
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Deposited on : 2001-09-06

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
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : rb-20027457
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027457

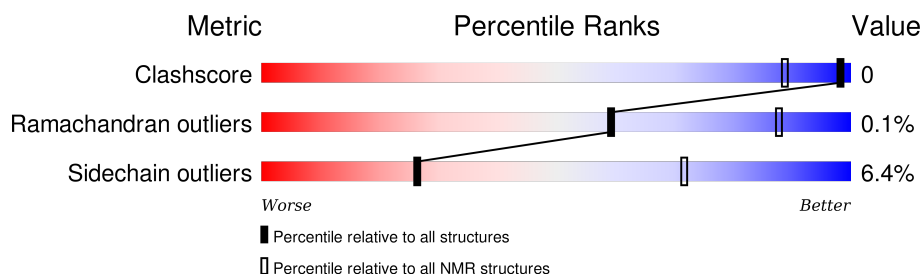
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 93%.

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	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

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	94	<div> <div></div> <div>80%</div> <div>5%</div> <div>15%</div> </div>

2 Ensemble composition and analysis

This entry contains 35 models. Model 2 is the overall representative, medoid model (most similar to other models). The authors have identified model 9 as representative.

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:33, A:41-A:98 (80)	0.33	2

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 5 clusters and 4 single-model clusters were found.

Cluster number	Models
1	1, 8, 9, 10, 15, 16, 21, 27, 28, 34
2	7, 13, 17, 20, 23, 26, 29, 31, 33
3	2, 3, 6, 11, 14, 22, 35
4	4, 5, 25
5	19, 32
Single-model clusters	12; 18; 24; 30

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 1393 atoms, of which 706 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called PROTEIN TYROSINE PHOSPHATASE.

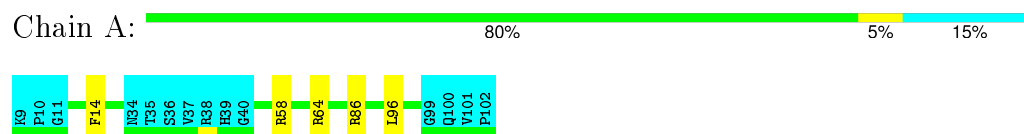
Mol	Chain	Residues	Atoms					Trace
1	A	94	Total	C	H	N	O	0
			1393	426	706	127	134	

4 Residue-property plots [i](#)

4.1 Average score per residue in the NMR ensemble

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

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE

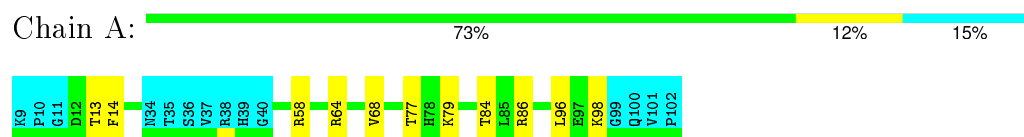


4.2 Scores per residue for each member of the ensemble

Colouring as in section [4.1](#) above.

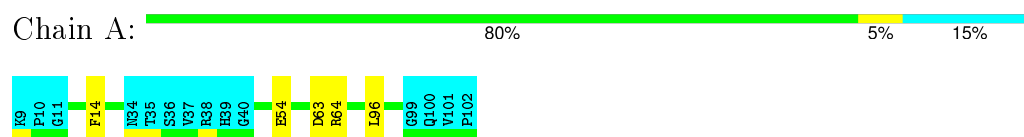
4.2.1 Score per residue for model 1

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



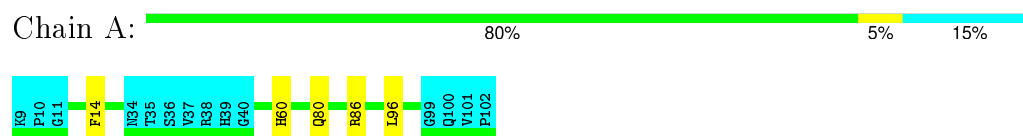
4.2.2 Score per residue for model 2 (medoid)

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



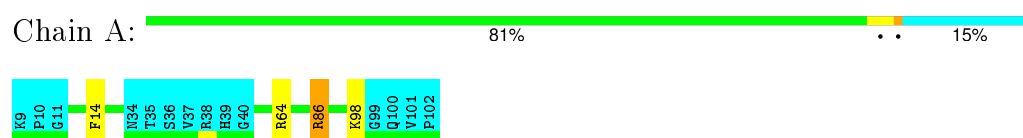
4.2.3 Score per residue for model 3

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



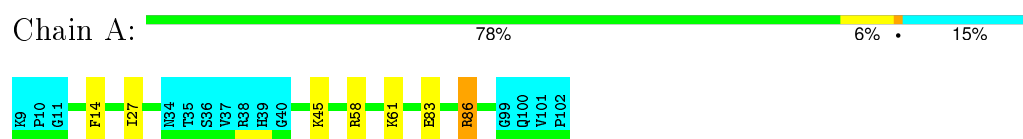
4.2.4 Score per residue for model 4

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



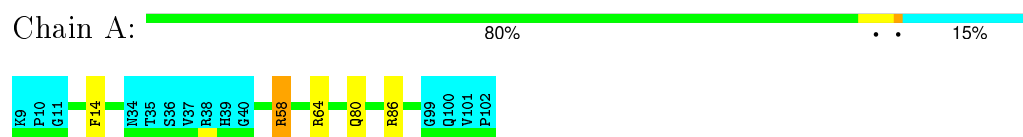
4.2.5 Score per residue for model 5

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



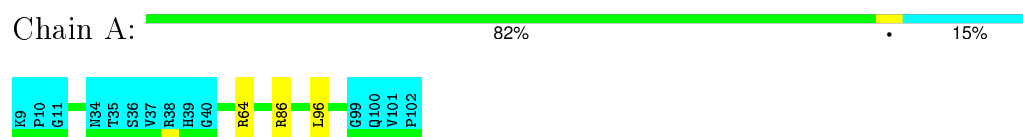
4.2.6 Score per residue for model 6

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



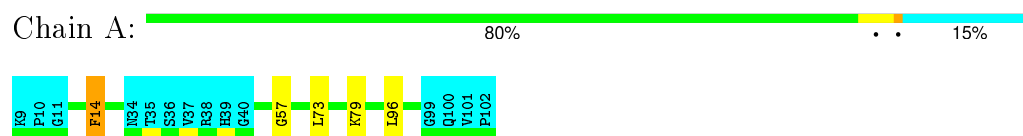
4.2.7 Score per residue for model 7

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



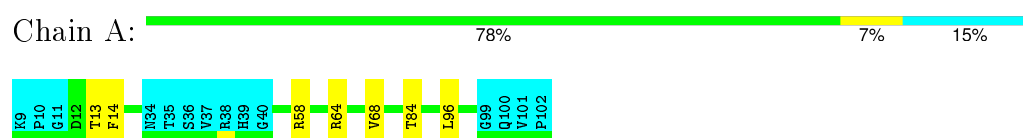
4.2.8 Score per residue for model 8

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



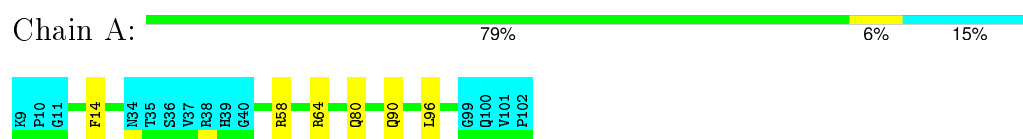
4.2.9 Score per residue for model 9

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



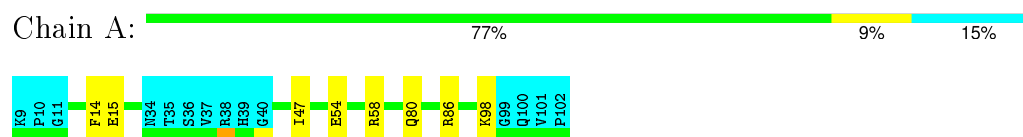
4.2.10 Score per residue for model 10

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



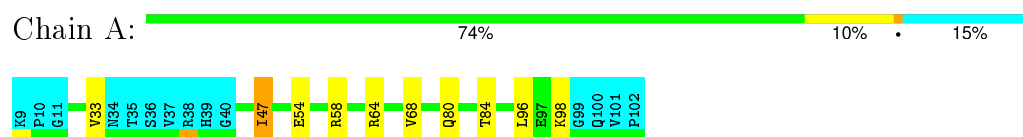
4.2.11 Score per residue for model 11

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



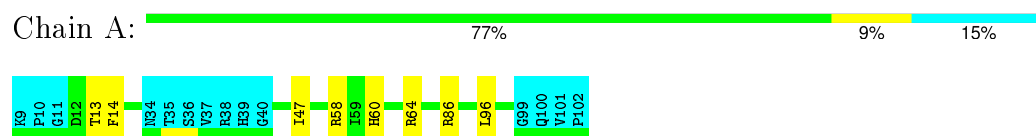
4.2.12 Score per residue for model 12

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



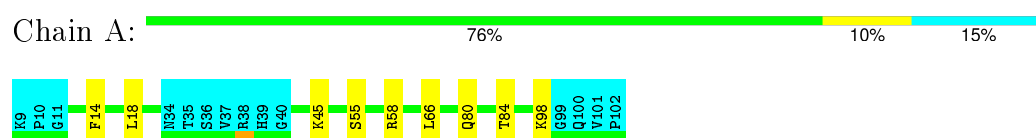
4.2.13 Score per residue for model 13

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



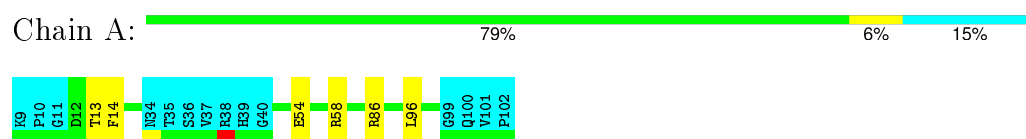
4.2.14 Score per residue for model 14

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



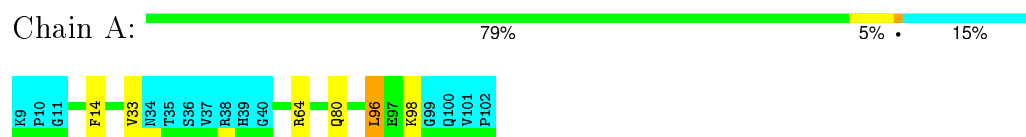
4.2.15 Score per residue for model 15

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



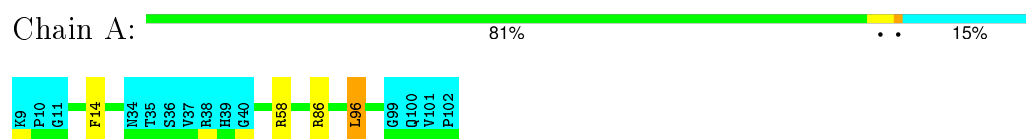
4.2.16 Score per residue for model 16

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



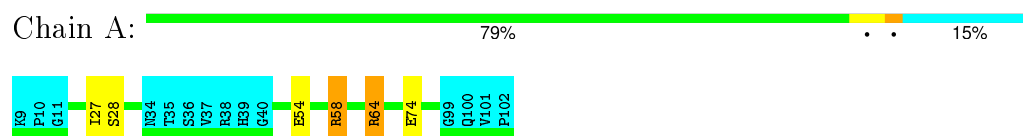
4.2.17 Score per residue for model 17

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



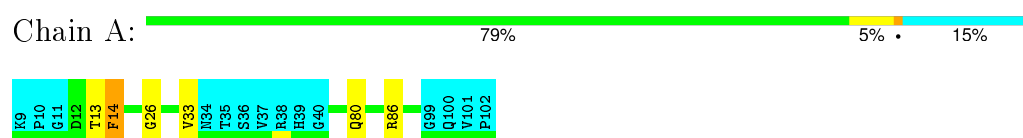
4.2.18 Score per residue for model 18

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



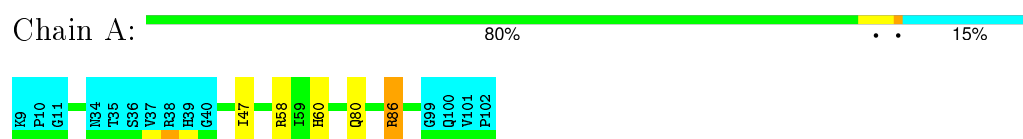
4.2.19 Score per residue for model 19

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



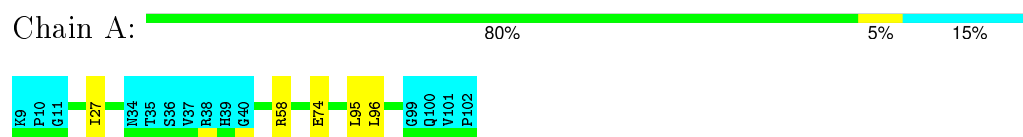
4.2.20 Score per residue for model 20

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



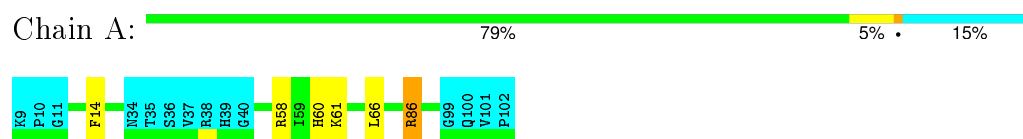
4.2.21 Score per residue for model 21

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



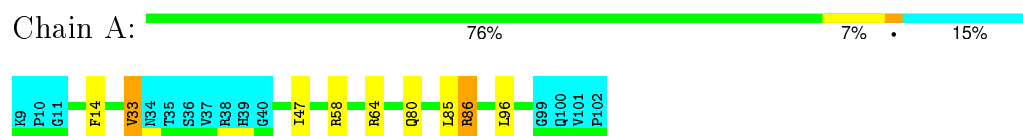
4.2.22 Score per residue for model 22

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



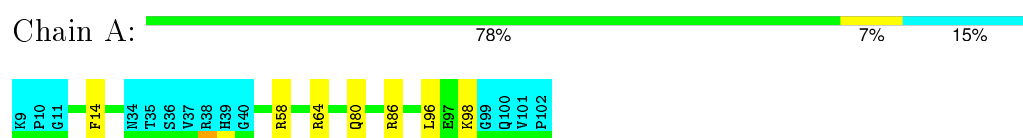
4.2.23 Score per residue for model 23

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



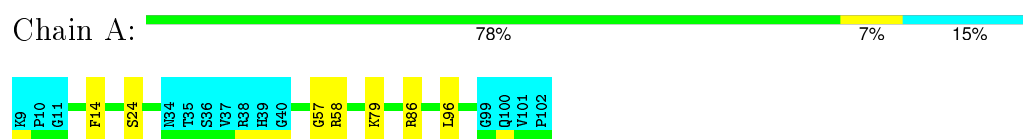
4.2.24 Score per residue for model 24

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



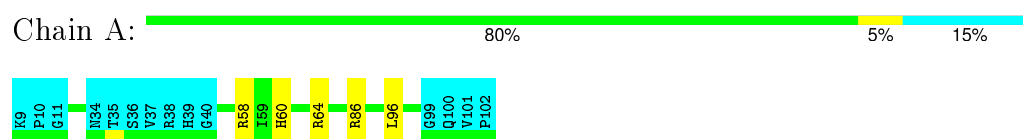
4.2.25 Score per residue for model 25

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



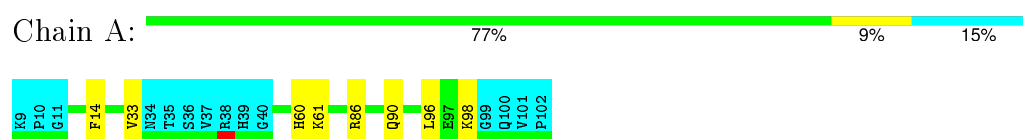
4.2.26 Score per residue for model 26

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



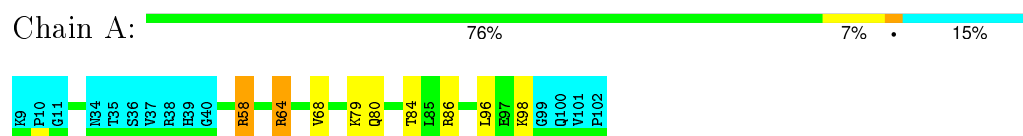
4.2.27 Score per residue for model 27

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



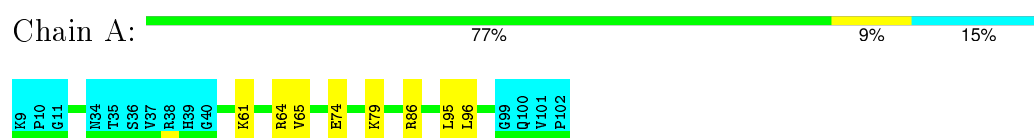
4.2.28 Score per residue for model 28

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



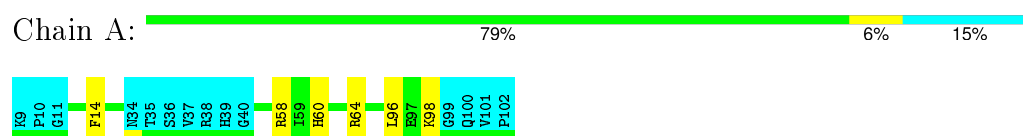
4.2.29 Score per residue for model 29

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



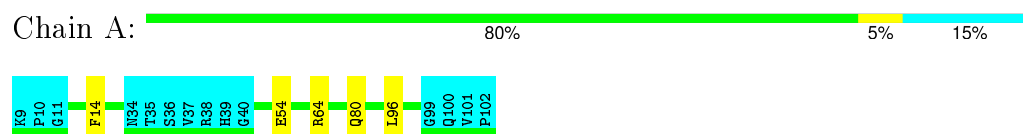
4.2.30 Score per residue for model 30

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



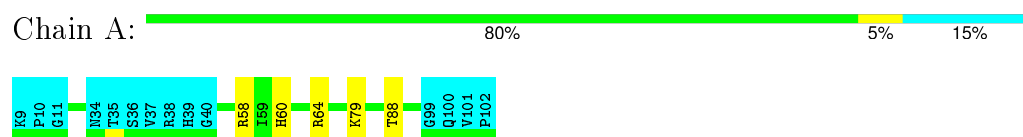
4.2.31 Score per residue for model 31

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



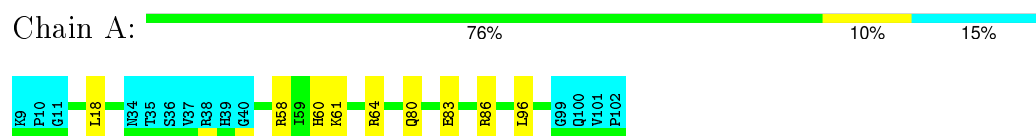
4.2.32 Score per residue for model 32

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



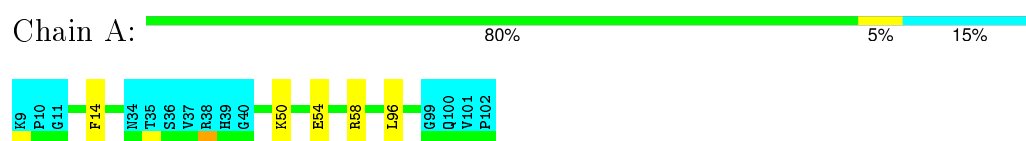
4.2.33 Score per residue for model 33

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



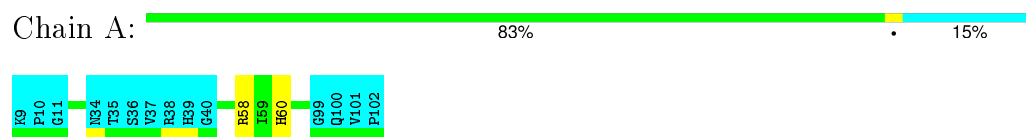
4.2.34 Score per residue for model 34

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



4.2.35 Score per residue for model 35

- Molecule 1: PROTEIN TYROSINE PHOSPHATASE



5 Refinement protocol and experimental data overview

The models were refined using the following method: *SIMULATED ANNEALING*, *RESTRAINED MOLECULAR DYNAMICS*.

Of the 200 calculated structures, 35 were deposited, based on the following criterion: *LOWEST ENERGY*.

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

Software name	Classification	Version
X-PLOR	refinement	
XPLOR	structure solution	

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 6060, BMRB entry 5131, BMRB entry 6092, 1
Number of chemical shift lists	
Total number of shifts	
Number of shifts mapped to atoms	
Number of unparsed shifts	
Number of shifts with mapping errors	
Number of shifts with mapping warnings	
Assignment completeness (well-defined parts)	

No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality

6.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	1.01±0.00	0±0/592 (0.0±0.0%)	1.08±0.03	2±1/800 (0.2±0.1%)
All	All	1.01	0/20720 (0.0%)	1.08	59/28000 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	0.6±0.6
All	All	0	21

There are no bond-length outliers.

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	14	PHE	CB-CG-CD2	-8.51	114.84	120.80	19	2
1	A	64	ARG	NE-CZ-NH1	8.19	124.39	120.30	2	15
1	A	58	ARG	NE-CZ-NH1	8.14	124.37	120.30	23	17
1	A	14	PHE	CB-CG-CD1	7.84	126.29	120.80	19	2
1	A	86	ARG	NE-CZ-NH1	6.69	123.64	120.30	11	13
1	A	58	ARG	NE-CZ-NH2	-6.29	117.16	120.30	23	5
1	A	64	ARG	NE-CZ-NH2	-5.78	117.41	120.30	2	4
1	A	86	ARG	NE-CZ-NH2	-5.33	117.63	120.30	11	1

There are no chirality outliers.

All unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	58	ARG	Sidechain	10
1	A	64	ARG	Sidechain	4
1	A	86	ARG	Sidechain	3
1	A	57	GLY	Peptide	2
1	A	27	ILE	Peptide	1
1	A	14	PHE	Sidechain	1

6.2 Too-close contacts ⓘ

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	586	606	606	1±1
All	All	20510	21210	21210	21

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 1.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:68:VAL:HG11	1:A:84:THR:HG21	0.80	1.53	12	4
1:A:60:HIS:CG	1:A:61:LYS:H	0.60	2.13	27	3
1:A:14:PHE:CZ	1:A:96:LEU:HD22	0.57	2.35	17	1
1:A:68:VAL:CG1	1:A:84:THR:HG21	0.53	2.31	12	2
1:A:14:PHE:CE1	1:A:96:LEU:HD21	0.52	2.39	16	1
1:A:60:HIS:CG	1:A:61:LYS:N	0.48	2.82	27	3
1:A:68:VAL:HG11	1:A:84:THR:CG2	0.47	2.38	28	2
1:A:14:PHE:CE2	1:A:96:LEU:HD11	0.47	2.45	10	1
1:A:14:PHE:CZ	1:A:96:LEU:HD21	0.45	2.46	16	1
1:A:14:PHE:CZ	1:A:96:LEU:HD11	0.42	2.50	25	1
1:A:47:ILE:HG23	1:A:54:GLU:HG2	0.42	1.92	12	2

6.3 Torsion angles ⓘ

6.3.1 Protein backbone ⓘ

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	80/94 (85%)	74±2 (93±2%)	6±2 (7±2%)	0±0 (0±0%)	59	88
All	All	2800/3290 (85%)	2603 (93%)	193 (7%)	4 (0%)	59	88

All 3 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	33	VAL	2
1	A	26	GLY	1
1	A	74	GLU	1

6.3.2 Protein sidechains ⓘ

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	62/73 (85%)	58±2 (94±3%)	4±2 (6±3%)	26	72
All	All	2170/2555 (85%)	2031 (94%)	139 (6%)	26	72

All 34 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	96	LEU	22
1	A	14	PHE	19
1	A	80	GLN	14
1	A	86	ARG	10
1	A	98	LYS	10
1	A	60	HIS	7
1	A	79	LYS	6
1	A	54	GLU	5
1	A	13	THR	5
1	A	33	VAL	4
1	A	47	ILE	4
1	A	66	LEU	2
1	A	64	ARG	2

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Mol	Chain	Res	Type	Models (Total)
1	A	18	LEU	2
1	A	27	ILE	2
1	A	61	LYS	2
1	A	83	GLU	2
1	A	45	LYS	2
1	A	90	GLN	2
1	A	74	GLU	2
1	A	95	LEU	2
1	A	77	THR	1
1	A	50	LYS	1
1	A	65	VAL	1
1	A	63	ASP	1
1	A	85	LEU	1
1	A	55	SER	1
1	A	73	LEU	1
1	A	24	SER	1
1	A	28	SER	1
1	A	88	THR	1
1	A	15	GLU	1
1	A	58	ARG	1
1	A	84	THR	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 ⓘ

There are no carbohydrates in this entry.

6.6 Ligand geometry ⓘ

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 93% for the well-defined parts and 92% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 5131

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	1073
Number of shifts mapped to atoms	1073
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	83	-0.18 ± 0.15	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	74	0.05 ± 0.12	None needed (< 0.5 ppm)
$^{13}\text{C}'$	86	-0.18 ± 0.07	None needed (< 0.5 ppm)
^{15}N	88	0.26 ± 0.28	None needed (< 0.5 ppm)

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 90%, i.e. 822 atoms were assigned a chemical shift out of a possible 914. 14 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	385/398 (97%)	158/159 (99%)	149/160 (93%)	78/79 (99%)
Sidechain	403/475 (85%)	245/271 (90%)	152/185 (82%)	6/19 (32%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	34/41 (83%)	14/21 (67%)	14/14 (100%)	6/6 (100%)
Overall	822/914 (90%)	417/451 (92%)	315/359 (88%)	90/104 (87%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 86%, i.e. 925 atoms were assigned a chemical shift out of a possible 1076. 15 out of 20 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	439/464 (95%)	182/185 (98%)	169/188 (90%)	88/91 (97%)
Sidechain	446/563 (79%)	273/324 (84%)	167/214 (78%)	6/25 (24%)
Aromatic	40/49 (82%)	16/25 (64%)	16/16 (100%)	8/8 (100%)
Overall	925/1076 (86%)	471/534 (88%)	352/418 (84%)	102/124 (82%)

7.1.4 Statistically unusual chemical shifts [i](#)

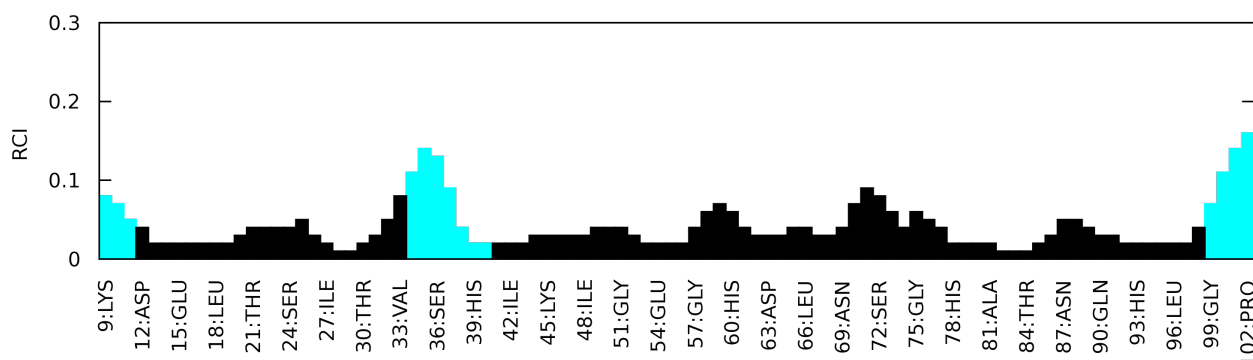
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	14	PHE	CE1	118.72	137.92 – 123.42	-8.2

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:



7.2 Chemical shift list 2

File name: BMRB entry 6060

Chemical shift list name: *assigned_chem_shift_list_1*

7.2.1 Bookkeeping [i](#)

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	1010
Number of shifts mapped to atoms	1010
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

7.2.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	90	-0.26 ± 0.26	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	79	0.14 ± 0.16	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	—
^{15}N	84	0.55 ± 0.43	None needed (imprecise)

7.2.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 81%, i.e. 743 atoms were assigned a chemical shift out of a possible 914. 17 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	311/398 (78%)	158/159 (99%)	77/160 (48%)	76/79 (96%)
Sidechain	405/475 (85%)	246/271 (91%)	158/185 (85%)	1/19 (5%)
Aromatic	27/41 (66%)	16/21 (76%)	6/14 (43%)	5/6 (83%)
Overall	743/914 (81%)	420/451 (93%)	241/359 (67%)	82/104 (79%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 80%, i.e. 863 atoms were assigned a chemical shift out of a possible 1076. 19 out of 20 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	356/464 (77%)	182/185 (98%)	90/188 (48%)	84/91 (92%)
Sidechain	473/563 (84%)	290/324 (90%)	182/214 (85%)	1/25 (4%)
Aromatic	34/49 (69%)	19/25 (76%)	8/16 (50%)	7/8 (88%)
Overall	863/1076 (80%)	491/534 (92%)	280/418 (67%)	92/124 (74%)

7.2.4 Statistically unusual chemical shifts [i](#)

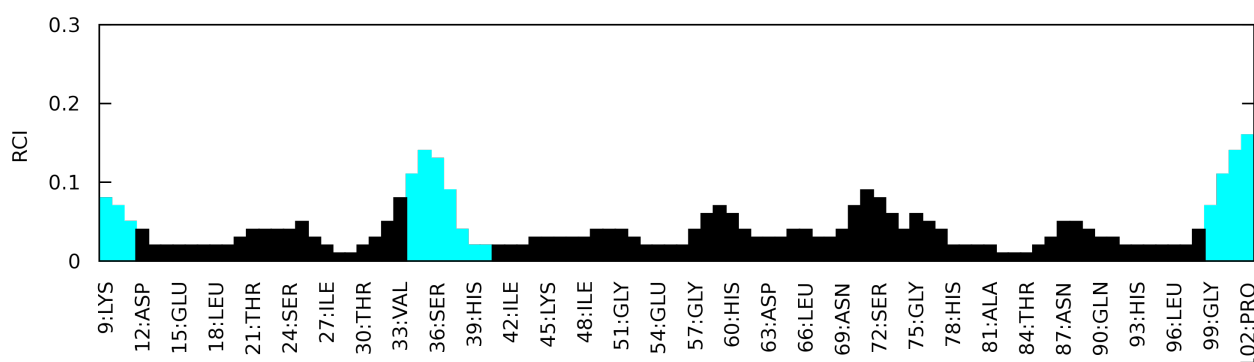
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	86	ARG	NE	121.69	92.63 – 76.73	23.3

7.2.5 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:



7.3 Chemical shift list 3

File name: BMRB entry 6091

Chemical shift list name: *assigned_chem_shift_list_1*

7.3.1 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	709
Number of shifts mapped to atoms	709
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	2

7.3.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	93	-0.65 ± 0.17	Should be applied
$^{13}\text{C}_\beta$	0	—	—
$^{13}\text{C}'$	0	—	—
^{15}N	89	0.49 ± 0.22	None needed (< 0.5 ppm)

7.3.3 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 57%, i.e. 523 atoms were assigned a chemical shift out of a possible 911. 0 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	308/398 (77%)	150/159 (94%)	80/160 (50%)	78/79 (99%)
Sidechain	206/475 (43%)	200/271 (74%)	0/185 (0%)	6/19 (32%)
Aromatic	9/38 (24%)	9/21 (43%)	0/14 (0%)	0/3 (0%)
Overall	523/911 (57%)	359/451 (80%)	80/359 (22%)	84/101 (83%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 55%, i.e. 594 atoms were assigned a chemical shift out of a possible 1072. 0 out of 20 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	352/464 (76%)	170/185 (92%)	93/188 (49%)	89/91 (98%)
Sidechain	233/563 (41%)	225/324 (69%)	0/214 (0%)	8/25 (32%)
Aromatic	9/45 (20%)	9/25 (36%)	0/16 (0%)	0/4 (0%)
Overall	594/1072 (55%)	404/534 (76%)	93/418 (22%)	97/120 (81%)

7.3.4 Statistically unusual chemical shifts [i](#)

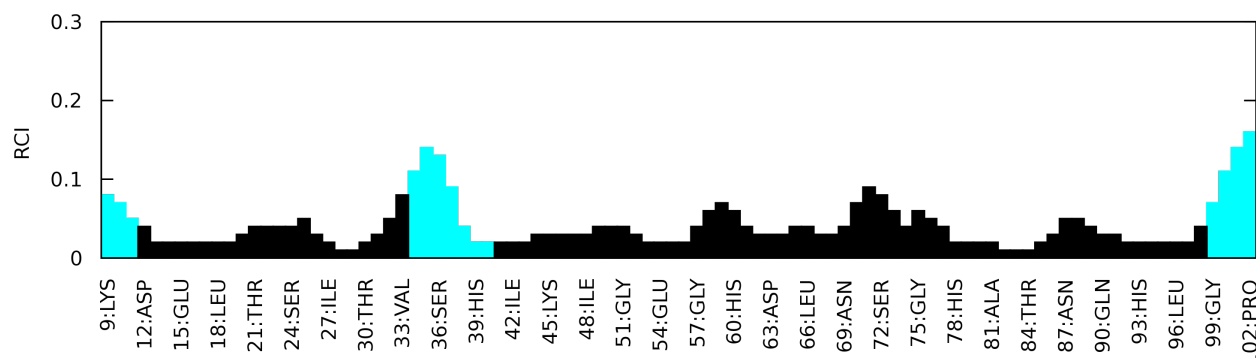
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, <i>ppm</i>	Expected range, <i>ppm</i>	Z-score
1	A	58	ARG	NE	123.18	92.63 – 76.73	24.2
1	A	64	ARG	NE	121.32	92.63 – 76.73	23.0

7.3.5 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:



7.4 Chemical shift list 4

File name: BMRB entry 6092

Chemical shift list name: *assigned_chem_shift_list_1*

7.4.1 Bookkeeping [i](#)

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	1005
Number of shifts mapped to atoms	1005
Number of unparsed shifts	0

Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	3

7.4.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	91	-0.30 ± 0.13	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	76	-0.04 ± 0.22	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	—
^{15}N	89	0.47 ± 0.22	None needed (< 0.5 ppm)

7.4.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 82%, i.e. 751 atoms were assigned a chemical shift out of a possible 914. 16 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	315/398 (79%)	158/159 (99%)	79/160 (49%)	78/79 (99%)
Sidechain	409/475 (86%)	247/271 (91%)	156/185 (84%)	6/19 (32%)
Aromatic	27/41 (66%)	15/21 (71%)	6/14 (43%)	6/6 (100%)
Overall	751/914 (82%)	420/451 (93%)	241/359 (67%)	90/104 (87%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 80%, i.e. 856 atoms were assigned a chemical shift out of a possible 1076. 17 out of 20 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	362/464 (78%)	182/185 (98%)	91/188 (48%)	89/91 (98%)
Sidechain	461/563 (82%)	280/324 (86%)	173/214 (81%)	8/25 (32%)
Aromatic	33/49 (67%)	17/25 (68%)	8/16 (50%)	8/8 (100%)
Overall	856/1076 (80%)	479/534 (90%)	272/418 (65%)	105/124 (85%)

7.4.4 Statistically unusual chemical shifts [i](#)

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, <i>ppm</i>	Expected range, <i>ppm</i>	Z-score
1	A	86	ARG	NE	123.94	92.63 – 76.73	24.7
1	A	58	ARG	NE	116.77	92.63 – 76.73	20.2
1	A	74	GLU	CG	29.09	42.24 – 29.94	-5.7

7.4.5 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:

