Full wwPDB NMR Structure Validation Report (i)

Mar 7, 2022 – 06:35 AM EST

PDB ID	:	2YUN
Title	:	Solution structure of the SH3 domain of human Nostrin
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Deposited on	:	2007-04-06

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
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:

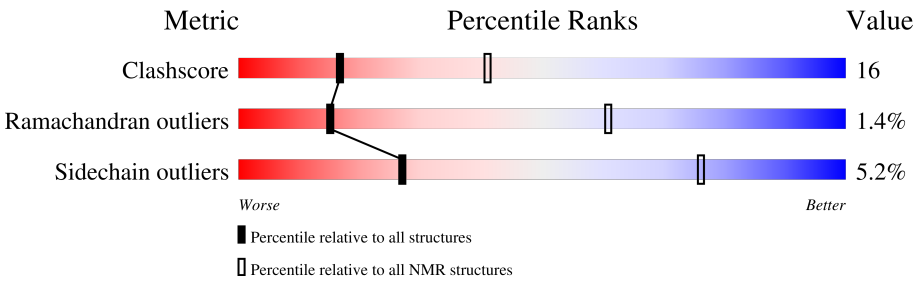
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.27
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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	79	48% 24% 28%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 3 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:7-A:63 (57)	0.19	3

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

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

The models can be grouped into 4 clusters and 2 single-model clusters were found.

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

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 1160 atoms, of which 567 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Nostrin.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	79	1160	369	567	104	119	1	0

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP Q8IVI9
A	2	SER	-	expression tag	UNP Q8IVI9
A	3	SER	-	expression tag	UNP Q8IVI9
A	4	GLY	-	expression tag	UNP Q8IVI9
A	5	SER	-	expression tag	UNP Q8IVI9
A	6	SER	-	expression tag	UNP Q8IVI9
A	7	GLY	-	expression tag	UNP Q8IVI9
A	74	SER	-	expression tag	UNP Q8IVI9
A	75	GLY	-	expression tag	UNP Q8IVI9
A	76	PRO	-	expression tag	UNP Q8IVI9
A	77	SER	-	expression tag	UNP Q8IVI9
A	78	SER	-	expression tag	UNP Q8IVI9
A	79	GLY	-	expression tag	UNP Q8IVI9

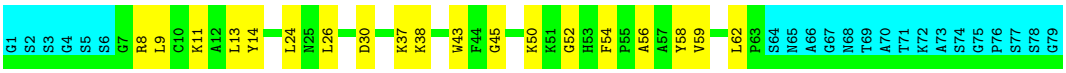
4 Residue-property plots (i)

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: Nostrin

Chain A: 48% 24% 28%



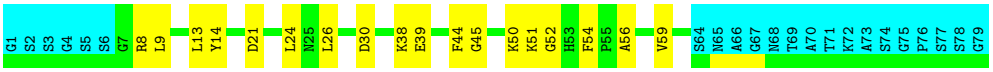
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

• Molecule 1: Nostrin

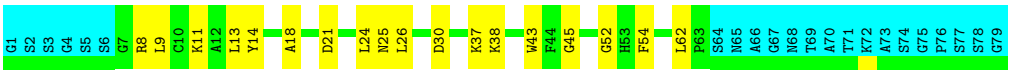
Chain A: 49% 23% 28%



4.2.2 Score per residue for model 2

• Molecule 1: Nostrin

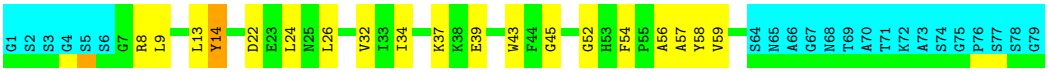
Chain A: 49% 23% 28%



4.2.3 Score per residue for model 3 (medoid)

- Molecule 1: Nostrin

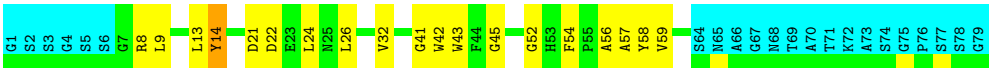
Chain A: 48% 23% • 28%



4.2.4 Score per residue for model 4

- Molecule 1: Nostrin

Chain A: 48% 23% • 28%



4.2.5 Score per residue for model 5

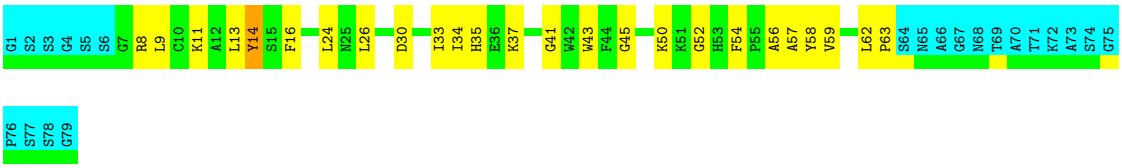
- Molecule 1: Nostrin

Chain A: 44% 28% 28%

4.2.6 Score per residue for model 6

- Molecule 1: Nostrin

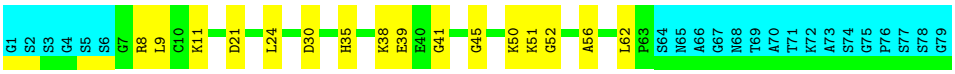
Chain A: 41% 30% • 28%



4.2.7 Score per residue for model 7

- Molecule 1: Nostrin

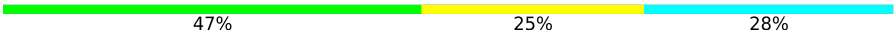
Chain A: 52% 20% 28%

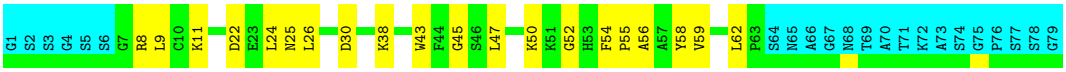


4.2.8 Score per residue for model 8

- Molecule 1: Nostrin

Chain A:

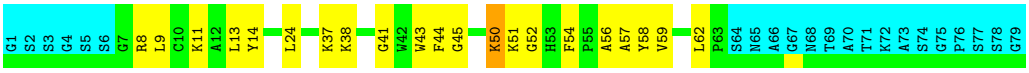




4.2.9 Score per residue for model 9

- Molecule 1: Nostrin

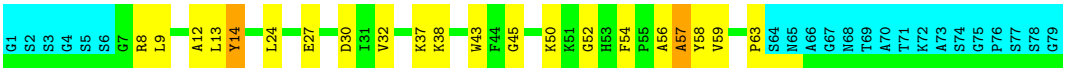
Chain A:



4.2.10 Score per residue for model 10

- Molecule 1: Nostrin

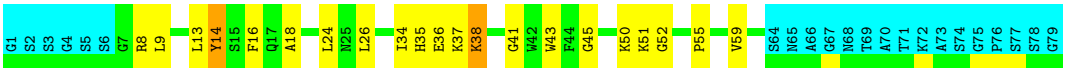
Chain A:



4.2.11 Score per residue for model 11

- Molecule 1: Nostrin

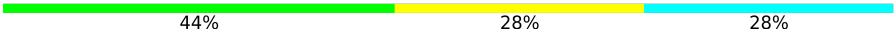
Chain A:

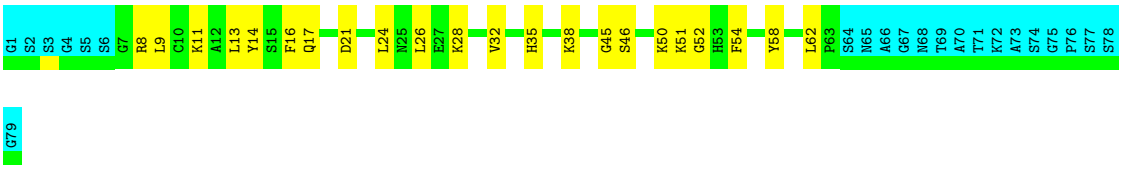


4.2.12 Score per residue for model 12

- Molecule 1: Nostrin

Chain A:

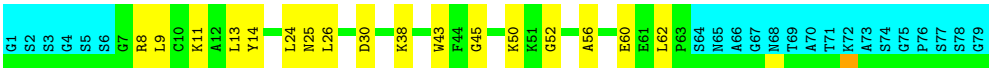




4.2.13 Score per residue for model 13

- Molecule 1: Nostrin

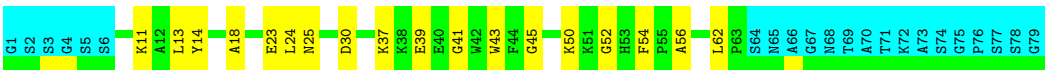
Chain A: 51% 22% 28%



4.2.14 Score per residue for model 14

- Molecule 1: Nostrin

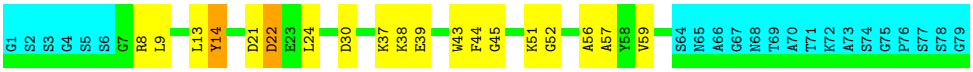
Chain A: 49% 23% 28%



4.2.15 Score per residue for model 15

- Molecule 1: Nostrin

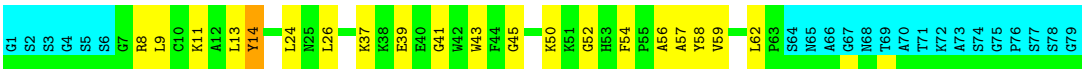
Chain A: 48% 22% • 28%



4.2.16 Score per residue for model 16

- Molecule 1: Nostrin

Chain A: 47% 24% • 28%



4.2.17 Score per residue for model 17

- Molecule 1: Nostrin

Chain A: 49% 23% 28%

G1 S2 S3 G4 S5 S6 L9 C10 K11 A12 L13 Y14 L24 N25 L26 I31 I34 K37 K38 G45 G52 H53 F54 P55 A56 V59 L62 P63 S64 N65 A66 G67 N68 T69 A70 T71 K72 A73 S74 G75 P76 S77 S78 G79

4.2.18 Score per residue for model 18

- Molecule 1: Nostrin

Chain A: 47% 25% 28%

G1 S2 S3 G4 S5 S6 G7 R8 L9 C10 K11 A12 L13 Y14 S15 F16 D21 L24 N25 L26 K37 K38 W42 W43 F44 G45 K50 K51 G52 L62 P63 S64 N65 A66 G67 N68 T69 A70 T71 K72 A73 S74 G75 P76 S77 S78 G79

4.2.19 Score per residue for model 19

- Molecule 1: Nostrin

Chain A: 41% 30% • 28%

G1 S2 S3 G4 S5 S6 G7 R8 L9 L13 Y14 L24 N25 L26 D30 I34 H35 E36 K37 K38 E39 E40 G41 W42 W43 F44 G45 K50 K51 G52 H53 F54 P55 A56 A57 Y58 V59 S64 N65 A66 G67 N68 T69 A70 T71 K72 A73 S74 G75 P76 S77 S78 G79

4.2.20 Score per residue for model 20

- Molecule 1: Nostrin

Chain A: 41% 32% 28%

G1 S2 S3 G4 S5 S6 G7 R8 L9 C10 K11 A12 L13 Y14 D21 L24 N25 D30 I31 I34 K37 K38 E39 E40 G41 W42 W43 F44 G45 S46 K50 K51 G52 H53 F54 P55 A56 A57 Y58 V59 L62 P63 S64 N65 A66 G67 N68 T69 A70 T71 K72 A73 S74

G75 P76 S77 S78 G79

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 violations, 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	464	451	449	14±4
All	All	9280	9020	8980	284

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 16.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:24:LEU:HD12	1:A:52:GLY:N	0.85	1.85	15	20
1:A:24:LEU:HD11	1:A:50:LYS:O	0.78	1.78	7	15
1:A:41:GLY:O	1:A:56:ALA:HB2	0.73	1.83	6	6
1:A:11:LYS:HB2	1:A:62:LEU:HD11	0.71	1.60	20	8
1:A:26:LEU:HD11	1:A:54:PHE:HB2	0.69	1.64	4	10
1:A:34:ILE:HG21	1:A:37:LYS:HE2	0.66	1.67	6	2
1:A:8:ARG:C	1:A:9:LEU:HD22	0.61	2.16	1	18
1:A:11:LYS:CB	1:A:62:LEU:HD11	0.58	2.27	18	9
1:A:34:ILE:HG21	1:A:37:LYS:HE3	0.57	1.75	11	2
1:A:35:HIS:CE1	1:A:51:LYS:NZ	0.57	2.72	11	1
1:A:41:GLY:O	1:A:43:TRP:CD1	0.57	2.57	4	4
1:A:34:ILE:HG21	1:A:37:LYS:HD3	0.55	1.78	17	1
1:A:24:LEU:HD12	1:A:52:GLY:CA	0.54	2.32	15	2
1:A:24:LEU:HD23	1:A:25:ASN:N	0.54	2.18	14	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:43:TRP:HB2	1:A:54:PHE:CE1	0.51	2.41	9	10
1:A:24:LEU:CD1	1:A:45:GLY:O	0.50	2.60	13	20
1:A:41:GLY:O	1:A:56:ALA:CB	0.50	2.60	16	7
1:A:56:ALA:O	1:A:59:VAL:N	0.49	2.45	16	12
1:A:34:ILE:HG21	1:A:37:LYS:CE	0.49	2.37	19	1
1:A:16:PHE:CD1	1:A:55:PRO:HG3	0.49	2.43	11	1
1:A:13:LEU:HB2	1:A:14:TYR:CD1	0.49	2.42	16	18
1:A:13:LEU:HB2	1:A:14:TYR:CE1	0.49	2.42	16	11
1:A:43:TRP:CD1	1:A:56:ALA:HB2	0.49	2.42	13	1
1:A:44:PHE:CE2	1:A:51:LYS:HD2	0.49	2.43	1	5
1:A:25:ASN:OD1	1:A:26:LEU:N	0.49	2.46	13	1
1:A:56:ALA:O	1:A:58:TYR:N	0.49	2.46	16	9
1:A:9:LEU:HD22	1:A:9:LEU:N	0.48	2.24	2	5
1:A:24:LEU:N	1:A:53:HIS:O	0.47	2.47	17	1
1:A:37:LYS:HD2	1:A:43:TRP:CE2	0.47	2.45	16	9
1:A:35:HIS:O	1:A:36:GLU:CG	0.47	2.63	5	3
1:A:12:ALA:HB3	1:A:27:GLU:O	0.47	2.10	10	1
1:A:37:LYS:CD	1:A:43:TRP:CD2	0.47	2.98	6	1
1:A:11:LYS:HB3	1:A:62:LEU:HD11	0.47	1.86	13	2
1:A:35:HIS:CE1	1:A:51:LYS:HE2	0.46	2.45	12	2
1:A:13:LEU:O	1:A:28:LYS:CE	0.46	2.63	12	1
1:A:18:ALA:HB2	1:A:24:LEU:C	0.45	2.32	11	2
1:A:25:ASN:O	1:A:26:LEU:HD23	0.45	2.11	8	1
1:A:44:PHE:CZ	1:A:51:LYS:HD2	0.45	2.46	19	2
1:A:14:TYR:CD1	1:A:14:TYR:N	0.45	2.85	9	8
1:A:24:LEU:HD23	1:A:25:ASN:O	0.44	2.12	13	1
1:A:9:LEU:HD12	1:A:31:ILE:CG2	0.44	2.41	20	2
1:A:26:LEU:HD13	1:A:59:VAL:CG1	0.44	2.43	11	1
1:A:24:LEU:CD1	1:A:50:LYS:O	0.44	2.60	12	1
1:A:32:VAL:HG11	1:A:54:PHE:CE2	0.44	2.48	12	5
1:A:16:PHE:HB2	1:A:58:TYR:CD2	0.44	2.48	6	2
1:A:35:HIS:CE1	1:A:51:LYS:HZ1	0.43	2.30	11	1
1:A:37:LYS:HD3	1:A:43:TRP:CE3	0.43	2.47	6	1
1:A:8:ARG:O	1:A:34:ILE:N	0.43	2.52	20	2
1:A:36:GLU:OE1	1:A:38:LYS:CE	0.43	2.67	11	1
1:A:18:ALA:HB1	1:A:23:GLU:HB3	0.43	1.90	14	1
1:A:16:PHE:CB	1:A:26:LEU:HD12	0.42	2.44	18	1
1:A:56:ALA:C	1:A:58:TYR:N	0.42	2.73	6	8
1:A:16:PHE:CG	1:A:17:GLN:N	0.42	2.88	12	1
1:A:43:TRP:CB	1:A:54:PHE:CE1	0.42	3.02	8	1
1:A:37:LYS:CD	1:A:43:TRP:CE2	0.42	3.03	9	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:41:GLY:O	1:A:56:ALA:N	0.42	2.53	16	2
1:A:11:LYS:HB2	1:A:62:LEU:CD1	0.41	2.45	9	1
1:A:56:ALA:O	1:A:57:ALA:C	0.41	2.58	15	2
1:A:13:LEU:HD11	1:A:60:GLU:HB2	0.41	1.92	13	1
1:A:13:LEU:N	1:A:58:TYR:O	0.41	2.52	16	1
1:A:33:ILE:O	1:A:35:HIS:CD2	0.41	2.73	6	1
1:A:9:LEU:CD2	1:A:9:LEU:N	0.41	2.84	12	1
1:A:21:ASP:O	1:A:22:ASP:C	0.41	2.58	4	2
1:A:9:LEU:N	1:A:9:LEU:CD2	0.41	2.84	7	1
1:A:34:ILE:HG21	1:A:37:LYS:CD	0.41	2.44	17	1
1:A:56:ALA:O	1:A:59:VAL:O	0.40	2.39	20	2
1:A:37:LYS:HG2	1:A:43:TRP:CG	0.40	2.51	10	1
1:A:47:LEU:C	1:A:47:LEU:HD12	0.40	2.37	8	1
1:A:55:PRO:HB2	1:A:58:TYR:CD1	0.40	2.52	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	57/79 (72%)	50±2 (88±3%)	6±2 (11±3%)	1±1 (1±2%)	15	61
All	All	1140/1580 (72%)	1001 (88%)	123 (11%)	16 (1%)	15	61

All 4 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	57	ALA	7
1	A	39	GLU	6
1	A	42	TRP	2
1	A	38	LYS	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	47/61 (77%)	45±1 (95±2%)	2±1 (5±2%)	27	76
All	All	940/1220 (77%)	891 (95%)	49 (5%)	27	76

All 10 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	38	LYS	15
1	A	30	ASP	11
1	A	14	TYR	7
1	A	21	ASP	6
1	A	22	ASP	3
1	A	39	GLU	2
1	A	46	SER	2
1	A	23	GLU	1
1	A	28	LYS	1
1	A	50	LYS	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