



Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 05:12 PM BST

PDB ID : 1TRL
Title : NMR SOLUTION STRUCTURE OF THE C-TERMINAL FRAGMENT 255-316 OF THERMOLYSIN: A DIMER FORMED BY SUBUNITS HAVING THE NATIVE STRUCTURE
Authors : Rico, M.; Jimenez, M.A.; Gonzalez, C.; De Filippis, V.; Fontana, A.
Deposited on : 1994-09-02

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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

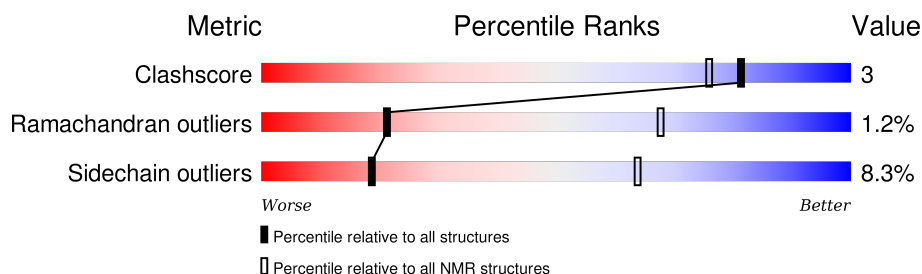
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment 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	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	62	
1	B	62	

2 Ensemble composition and analysis

This entry contains 8 models. Model 4 is the overall representative, medoid model (most similar to other models).

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:262-A:313, B:262-B:313 (104)	0.40	4

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 and 1 single-model cluster was found.

Cluster number	Models
1	1, 2, 3, 4, 6
2	7, 8
Single-model clusters	5

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1168 atoms, of which 232 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called THERMOLYSIN FRAGMENT 255 - 316.

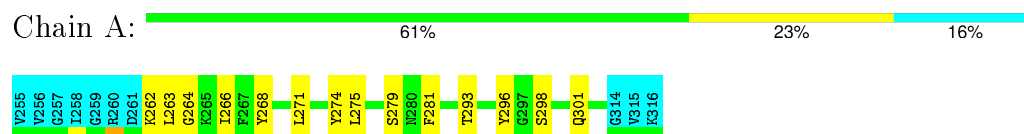
Mol	Chain	Residues	Atoms					Trace
1	A	62	Total	C	H	N	O	0
			584	296	116	81	91	
1	B	62	Total	C	H	N	O	0
			584	296	116	81	91	

4 Residue-property plots [i](#)

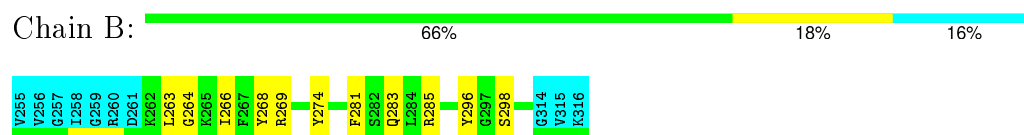
4.1 Average score per residue in the NMR ensemble

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: THERMOLYSIN FRAGMENT 255 - 316



- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

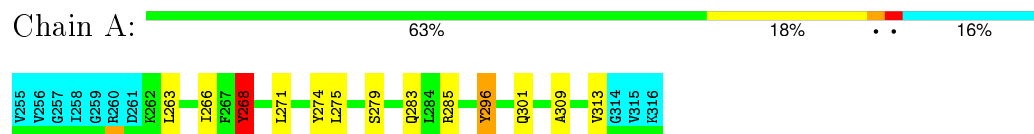


4.2 Scores per residue for each member of the ensemble

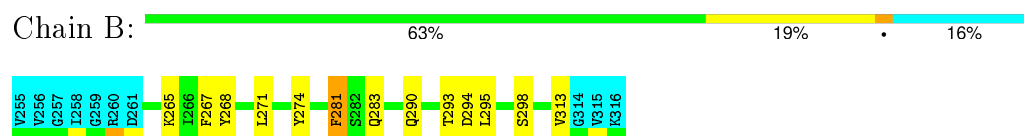
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

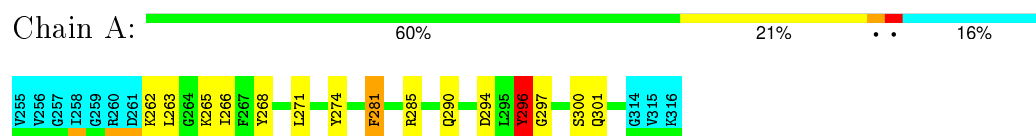


- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

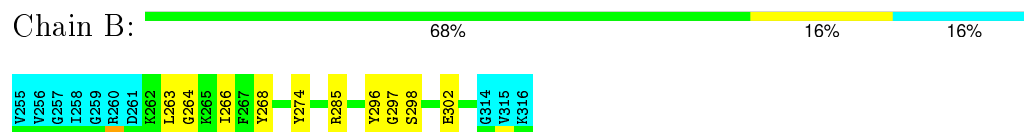


4.2.2 Score per residue for model 2

- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

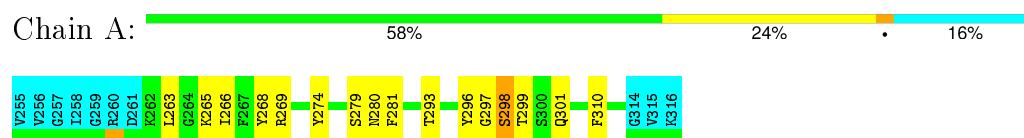


- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

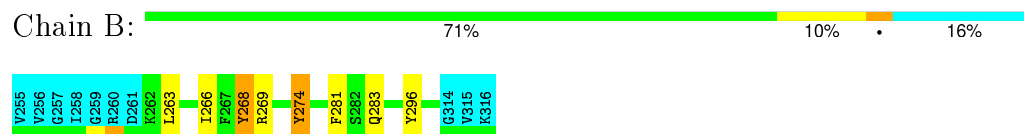


4.2.3 Score per residue for model 3

- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316



- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

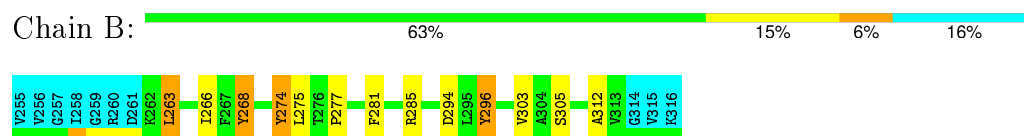


4.2.4 Score per residue for model 4 (medoid)

- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

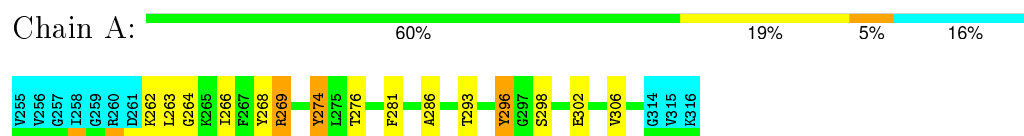


- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

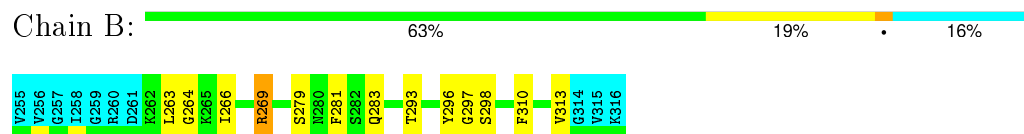


4.2.5 Score per residue for model 5

- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

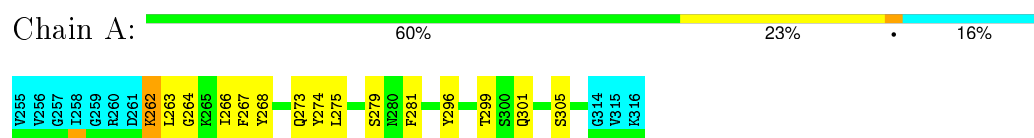


- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

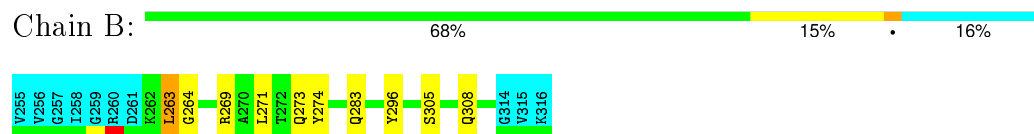


4.2.6 Score per residue for model 6

- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

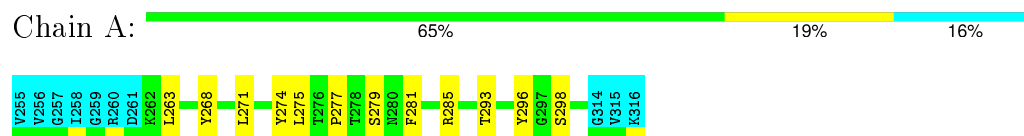


- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

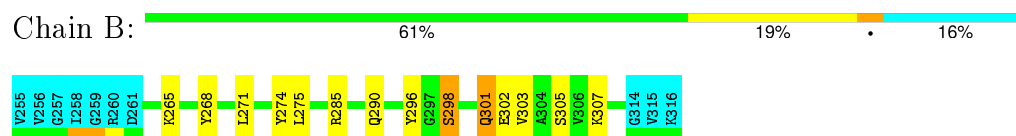


4.2.7 Score per residue for model 7

- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316



- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316



4.2.8 Score per residue for model 8

- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

Chain A:  61% 19% 16%



- Molecule 1: THERMOLYSIN FRAGMENT 255 - 316

Chain B:  68% 15% 16%



5 Refinement protocol and experimental data overview

Of the ? calculated structures, 8 were deposited, based on the following criterion: ?.

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

Software name	Classification	Version
GROMOS	refinement	

No chemical shift data was provided. 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	0.81±0.01	0±0/405 (0.0±0.0%)	1.26±0.10	3±2/550 (0.6±0.4%)
1	B	0.81±0.02	0±0/405 (0.0±0.0%)	1.24±0.08	3±2/550 (0.5±0.3%)
All	All	0.81	0/6480 (0.0%)	1.25	46/8800 (0.5%)

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	3.1±1.5
1	B	0.0±0.0	2.6±1.6
All	All	0	46

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	301	GLN	CG-CD-OE1	10.35	142.31	121.60	4	1
1	A	301	GLN	CG-CD-NE2	-8.56	96.14	116.70	4	1
1	A	274	TYR	CB-CG-CD2	-8.16	116.11	121.00	4	7
1	B	268	TYR	CB-CG-CD2	-7.54	116.48	121.00	3	1
1	A	285	ARG	NE-CZ-NH2	7.35	123.97	120.30	7	1
1	B	285	ARG	CD-NE-CZ	7.29	133.80	123.60	4	1
1	B	268	TYR	CB-CG-CD1	-7.28	116.63	121.00	4	1
1	B	269	ARG	NE-CZ-NH1	7.24	123.92	120.30	5	3
1	A	285	ARG	NE-CZ-NH1	6.87	123.73	120.30	1	2
1	A	296	TYR	CB-CG-CD2	-6.67	117.00	121.00	2	5
1	B	303	VAL	CG1-CB-CG2	-6.26	100.89	110.90	4	1
1	B	281	PHE	CB-CG-CD2	-6.24	116.43	120.80	5	3
1	B	296	TYR	CB-CG-CD2	-6.13	117.32	121.00	2	2

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	B	274	TYR	CB-CG-CD2	-6.08	117.35	121.00	6	4
1	A	267	PHE	CB-CG-CD1	-6.01	116.60	120.80	4	2
1	B	285	ARG	CG-CD-NE	-5.96	99.29	111.80	4	1
1	B	285	ARG	NE-CZ-NH1	5.89	123.25	120.30	8	1
1	A	268	TYR	CB-CG-CD2	-5.89	117.47	121.00	1	3
1	A	281	PHE	CB-CG-CD2	-5.87	116.69	120.80	2	2
1	A	296	TYR	CB-CG-CD1	-5.74	117.56	121.00	2	1
1	A	269	ARG	NE-CZ-NH1	5.42	123.01	120.30	5	1
1	B	267	PHE	CB-CG-CD1	-5.31	117.08	120.80	1	1
1	B	310	PHE	CB-CG-CD1	-5.28	117.10	120.80	8	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	B	296	TYR	Sidechain	5
1	A	296	TYR	Sidechain	5
1	A	268	TYR	Sidechain	5
1	B	274	TYR	Sidechain	5
1	B	268	TYR	Sidechain	4
1	B	281	PHE	Sidechain	3
1	A	281	PHE	Sidechain	3
1	A	310	PHE	Sidechain	2
1	A	274	TYR	Sidechain	2
1	A	313	VAL	Mainchain	1
1	B	312	ALA	Mainchain	1
1	A	278	THR	Mainchain	1
1	B	277	PRO	Mainchain	1
1	A	286	ALA	Mainchain	1
1	A	267	PHE	Sidechain	1
1	A	309	ALA	Mainchain	1
1	A	279	SER	Mainchain	1
1	A	306	VAL	Mainchain	1
1	B	310	PHE	Sidechain	1
1	A	308	GLN	Sidechain	1
1	B	313	VAL	Mainchain	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	398	96	399	4±2
1	B	398	96	399	3±1
All	All	6368	1536	6384	41

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 3.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:264:GLY:HA2	1:B:263:LEU:HB3	0.69	1.64	4	2
1:A:264:GLY:HA2	1:B:263:LEU:HB2	0.66	1.64	8	1
1:A:263:LEU:HA	1:A:266:ILE:HD12	0.62	1.70	6	6
1:B:269:ARG:CZ	1:B:273:GLN:HE22	0.59	2.09	6	1
1:A:263:LEU:HB2	1:B:264:GLY:HA2	0.58	1.76	8	1
1:A:293:THR:HG23	1:A:298:SER:HB3	0.57	1.76	5	3
1:B:293:THR:HG23	1:B:298:SER:HB3	0.54	1.80	5	2
1:A:263:LEU:HD22	1:B:264:GLY:HA2	0.52	1.80	6	3
1:B:265:LYS:HB3	1:B:295:LEU:HD13	0.50	1.82	1	1
1:A:300:SER:HB3	1:A:303:VAL:HG23	0.49	1.84	4	1
1:A:281:PHE:HB2	1:B:279:SER:HB2	0.49	1.84	5	1
1:B:263:LEU:HA	1:B:266:ILE:HD12	0.47	1.85	3	4
1:A:309:ALA:O	1:A:313:VAL:HG23	0.47	2.10	1	2
1:B:266:ILE:HD11	1:B:302:GLU:HG2	0.46	1.85	2	1
1:A:305:SER:HA	1:A:308:GLN:OE1	0.44	2.12	8	1
1:A:266:ILE:HG21	1:A:306:VAL:HG21	0.44	1.90	5	1
1:A:303:VAL:HG12	1:A:307:LYS:HE3	0.43	1.90	4	1
1:A:275:LEU:HB3	1:B:313:VAL:HG13	0.43	1.89	1	1
1:A:297:GLY:O	1:A:300:SER:HB2	0.43	2.13	2	1
1:B:271:LEU:HA	1:B:275:LEU:HB2	0.42	1.90	7	1
1:A:264:GLY:HA2	1:B:263:LEU:HD12	0.41	1.90	5	1
1:B:269:ARG:NH2	1:B:273:GLN:HE22	0.41	2.12	6	1
1:A:275:LEU:HG	1:A:279:SER:OG	0.41	2.15	6	1
1:A:293:THR:HG22	1:A:298:SER:HA	0.41	1.91	8	1
1:B:298:SER:HA	1:B:303:VAL:HG21	0.41	1.91	7	1
1:A:271:LEU:HA	1:A:275:LEU:HB2	0.41	1.93	4	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	52/62 (84%)	48±2 (92±3%)	3±2 (6±3%)	1±1 (2±2%)	16	59
1	B	52/62 (84%)	48±1 (91±2%)	4±1 (8±2%)	0±0 (1±1%)	31	76
All	All	832/992 (84%)	764 (92%)	58 (7%)	10 (1%)	21	68

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	299	THR	3
1	B	297	GLY	2
1	A	297	GLY	1
1	A	296	TYR	1
1	B	264	GLY	1
1	A	279	SER	1
1	A	298	SER	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	42/49 (86%)	38±1 (91±2%)	4±1 (9±2%)	18	63
1	B	42/49 (86%)	39±2 (92±4%)	3±2 (8±4%)	20	65
All	All	672/784 (86%)	616 (92%)	56 (8%)	19	64

All 36 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	271	LEU	5
1	B	283	GLN	4
1	A	301	GLN	3
1	A	265	LYS	3
1	A	279	SER	2
1	B	271	LEU	2
1	A	262	LYS	2
1	B	290	GLN	2
1	B	263	LEU	2
1	B	285	ARG	2
1	B	294	ASP	2
1	B	298	SER	2
1	A	275	LEU	2
1	B	308	GLN	1
1	A	294	ASP	1
1	B	269	ARG	1
1	B	302	GLU	1
1	A	280	ASN	1
1	A	302	GLU	1
1	B	273	GLN	1
1	B	301	GLN	1
1	A	276	THR	1
1	B	307	LYS	1
1	A	269	ARG	1
1	B	265	LYS	1
1	B	275	LEU	1
1	A	273	GLN	1
1	B	305	SER	1
1	A	263	LEU	1
1	A	283	GLN	1
1	A	268	TYR	1
1	A	293	THR	1
1	A	277	PRO	1
1	B	268	TYR	1
1	B	299	THR	1
1	A	290	GLN	1

6.3.3 RNA ⓘ

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 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

No chemical shift data were provided