



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:21 PM UTC

PDB ID : 5TUC / pdb_00005tuc
Title : Crystal Structure of the Sus TBC1D15 GAP Domain
Authors : Chen, Y.-N.; Wang, W.; Cheng, D.; Ge, Y.; Gu, X.; Zhou, X.E.; Ye, F.; Xu, H.E.; Lv, Z.
Deposited on : 2016-11-05
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray 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/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

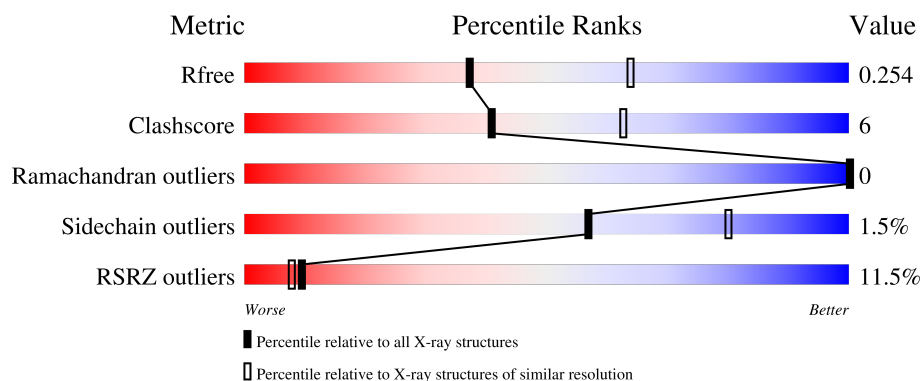
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	348	8% 77% 17% • 5%
1	B	348	13% 76% 14% • 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5559 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Sus TBC1D15 GAP Domain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	0	0
			2772	1771	466	510	25			
1	B	315	Total	C	N	O	S	0	0	0
			2633	1688	440	480	25			

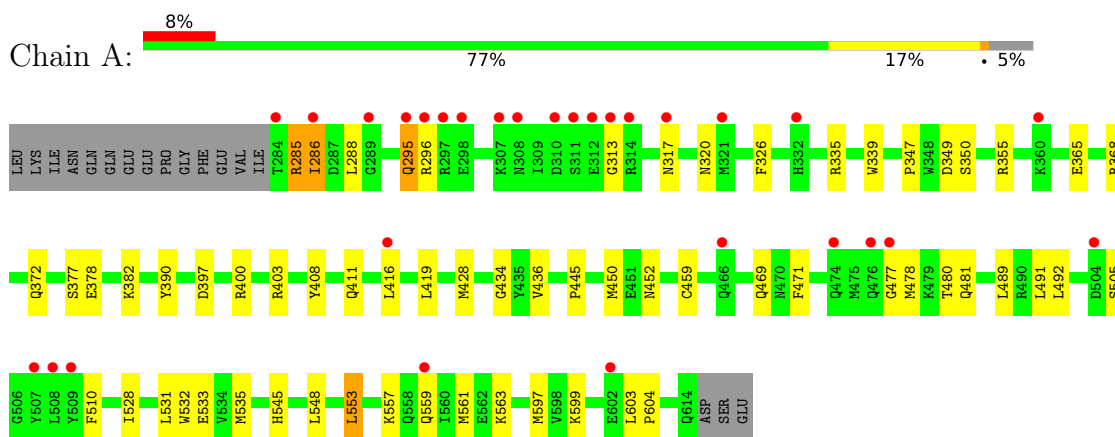
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	106	Total	O	0	0
			106	106		
2	B	48	Total	O	0	0
			48	48		

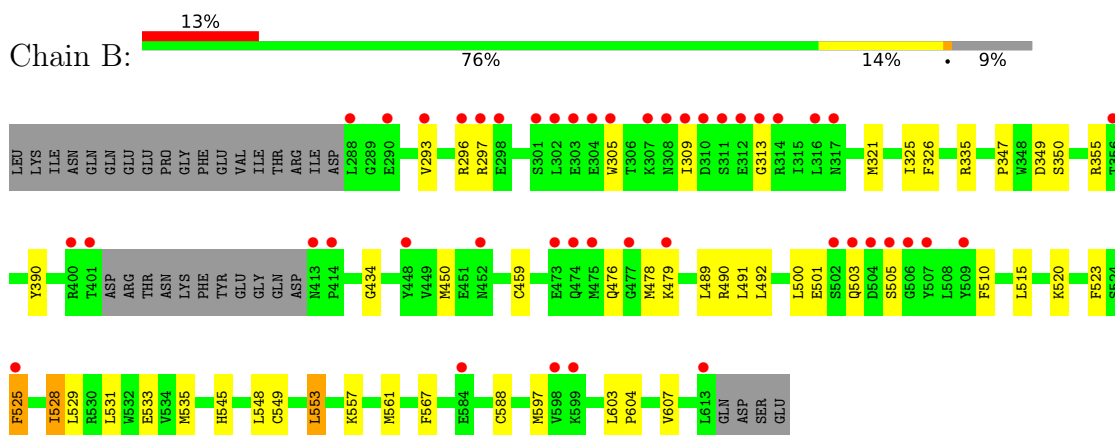
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Sus TBC1D15 GAP Domain



• Molecule 1: Sus TBC1D15 GAP Domain



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	139.00Å 139.00Å 175.24Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	87.62 – 2.50 87.62 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (87.62-2.50) 100.0 (87.62-2.50)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.214 , 0.253 0.217 , 0.254	Depositor DCC
R_{free} test set	2500 reflections (7.10%)	wwPDB-VP
Wilson B-factor (Å ²)	38.8	Xtriage
Anisotropy	0.024	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5559	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.76% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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 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	0.34	0/2830	0.76	1/3806 (0.0%)
1	B	0.28	0/2688	0.72	0/3614
All	All	0.31	0/5518	0.74	1/7420 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	478	MET	N-CA-C	-5.37	105.09	111.69

There are no chirality outliers.

There are no planarity outliers.

5.2 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 the 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 within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2772	0	2742	40	0
1	B	2633	0	2616	32	0
2	A	106	0	0	5	0
2	B	48	0	0	1	0
All	All	5559	0	5358	70	0

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 6.

All (70) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347:PRO:HG2	1:A:350:SER:HB3	1.60	0.84
1:A:285:ARG:HD2	1:A:286:ILE:HD12	1.63	0.79
1:B:347:PRO:HG2	1:B:350:SER:HB3	1.65	0.78
1:B:545:HIS:HA	1:B:548:LEU:HD12	1.69	0.74
1:A:296:ARG:NH1	1:A:326:PHE:O	2.21	0.74
1:B:603:LEU:HD12	1:B:604:PRO:HD2	1.70	0.74
1:B:503:GLN:HG2	1:B:567:PHE:HB2	1.69	0.73
1:B:505:SER:HB3	1:B:510:PHE:HE2	1.56	0.71
1:B:313:GLY:O	1:B:355:ARG:NH2	2.25	0.70
1:B:296:ARG:NH2	1:B:326:PHE:O	2.25	0.69
1:A:313:GLY:O	1:A:355:ARG:NH2	2.25	0.68
1:A:491:LEU:HD23	1:A:597:MET:HE2	1.78	0.65
1:B:491:LEU:HD23	1:B:597:MET:HE2	1.79	0.65
1:B:335:ARG:NE	1:B:533:GLU:OE2	2.24	0.63
1:A:545:HIS:HA	1:A:548:LEU:HD12	1.80	0.62
1:A:403:ARG:NH2	1:A:411:GLN:OE1	2.33	0.62
1:A:557:LYS:HG2	1:A:561:MET:HE2	1.84	0.60
1:A:489:LEU:HD11	1:A:553:LEU:HD23	1.84	0.59
1:A:603:LEU:HD12	1:A:604:PRO:HD2	1.85	0.58
1:B:296:ARG:HG3	1:B:588:CYS:SG	2.44	0.58
1:A:335:ARG:NE	1:A:533:GLU:OE2	2.29	0.58
1:A:295:GLN:HG2	1:B:297:ARG:HA	1.87	0.57
1:A:477:GLY:HA2	1:A:480:THR:HB	1.86	0.57
1:A:599:LYS:HD2	1:B:293:VAL:HG23	1.87	0.55
1:B:557:LYS:HG2	1:B:561:MET:HE2	1.88	0.54
1:A:397:ASP:OD1	1:A:400:ARG:NH1	2.42	0.53
1:A:469:GLN:NE2	2:A:702:HOH:O	2.40	0.53
1:B:525:PHE:O	1:B:529:LEU:HG	2.09	0.53
1:B:490:ARG:NH2	1:B:501:GLU:OE2	2.40	0.52
1:A:372:GLN:HB3	1:A:428:MET:HE2	1.92	0.52
1:A:339:TRP:HE1	1:A:533:GLU:HG2	1.75	0.51
1:A:403:ARG:NH1	1:A:416:LEU:HD11	2.25	0.51
1:A:445:PRO:HG2	1:A:532:TRP:CZ3	2.46	0.50
1:B:531:LEU:HG	1:B:535:MET:HE3	1.94	0.50
1:A:365:GLU:HG3	1:A:368:ARG:NH2	2.26	0.49
1:B:489:LEU:HD11	1:B:553:LEU:HD23	1.94	0.49
1:A:450:MET:HE3	1:A:459:CYS:SG	2.52	0.49
1:A:505:SER:HB3	1:A:510:PHE:HE2	1.77	0.48
1:A:477:GLY:O	1:A:481:GLN:NE2	2.47	0.48
1:B:478:MET:HE1	1:B:515:LEU:HB2	1.95	0.47
1:B:523:PHE:HB2	1:B:528:ILE:HG12	1.97	0.47
1:A:339:TRP:HE1	1:A:533:GLU:CG	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:378:GLU:HG2	1:A:382:LYS:HE3	1.98	0.46
1:A:436:VAL:HB	1:A:471:PHE:HB3	1.97	0.46
1:A:492:LEU:HD21	1:A:597:MET:HE3	1.98	0.46
1:A:599:LYS:HB3	1:A:599:LYS:HE2	1.61	0.45
1:A:349:ASP:OD1	1:A:349:ASP:N	2.45	0.45
1:B:500:LEU:HD13	1:B:510:PHE:HZ	1.81	0.45
1:B:549:CYS:O	1:B:553:LEU:HB2	2.17	0.45
1:B:491:LEU:HD22	1:B:607:VAL:HG21	1.99	0.45
1:A:563:LYS:NZ	2:A:708:HOH:O	2.49	0.45
1:A:390:TYR:CZ	1:A:434:GLY:HA3	2.52	0.45
1:A:531:LEU:HG	1:A:535:MET:HE3	1.98	0.45
1:B:305:TRP:CD1	1:B:309:ILE:HD11	2.52	0.44
1:A:285:ARG:HA	1:A:285:ARG:HD3	1.63	0.44
1:B:535:MET:HE2	1:B:548:LEU:HD13	2.01	0.43
1:A:355:ARG:NH1	2:A:709:HOH:O	2.51	0.43
1:A:559:GLN:HG2	2:A:735:HOH:O	2.18	0.43
1:B:450:MET:HE3	1:B:459:CYS:SG	2.59	0.43
1:A:377:SER:OG	2:A:701:HOH:O	2.19	0.43
1:B:349:ASP:OD1	1:B:349:ASP:N	2.46	0.42
1:B:520:LYS:HA	1:B:528:ILE:HG13	2.01	0.42
1:A:450:MET:C	1:A:452:ASN:H	2.26	0.42
1:B:321:MET:O	1:B:325:ILE:HG13	2.20	0.42
1:B:492:LEU:HD21	1:B:597:MET:HE3	2.01	0.42
1:B:476:GLN:HA	1:B:479:LYS:HB2	2.01	0.42
1:A:408:TYR:CE1	1:A:419:LEU:HG	2.54	0.41
1:B:390:TYR:CZ	1:B:434:GLY:HA3	2.55	0.41
1:A:317:ASN:HB3	1:A:320:ASN:HB3	2.01	0.41
1:B:520:LYS:HD2	2:B:716:HOH:O	2.21	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 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 X-ray entries followed by that with respect to entries of similar resolution.

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	329/348 (94%)	323 (98%)	6 (2%)	0	100	100
1	B	311/348 (89%)	304 (98%)	7 (2%)	0	100	100
All	All	640/696 (92%)	627 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 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 X-ray entries followed by that with respect to entries of similar resolution.

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	309/325 (95%)	303 (98%)	6 (2%)	50	76
1	B	294/325 (90%)	291 (99%)	3 (1%)	68	86
All	All	603/650 (93%)	594 (98%)	9 (2%)	57	80

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	285	ARG
1	A	286	ILE
1	A	288	LEU
1	A	295	GLN
1	A	528	ILE
1	A	553	LEU
1	B	525	PHE
1	B	528	ILE
1	B	553	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	317	ASN
1	A	323	GLN
1	A	481	GLN
1	B	317	ASN

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Mol	Chain	Res	Type
1	B	466	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	331/348 (95%)	0.43	29 (8%) 15 14	25, 39, 75, 139	0
1	B	315/348 (90%)	0.97	45 (14%) 6 5	31, 55, 101, 140	0
All	All	646/696 (92%)	0.69	74 (11%) 9 8	25, 47, 94, 140	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	288	LEU	6.0
1	B	401	THR	5.4
1	A	297	ARG	5.3
1	B	296	ARG	5.2
1	B	507	TYR	5.0
1	A	313	GLY	4.9
1	B	302	LEU	4.8
1	A	310	ASP	4.6
1	A	296	ARG	4.1
1	A	284	THR	4.0
1	B	297	ARG	3.9
1	B	413	ASN	3.8
1	B	313	GLY	3.6
1	A	312	GLU	3.5
1	B	502	SER	3.5
1	A	314	ARG	3.4
1	A	295	GLN	3.4
1	B	312	GLU	3.3
1	B	506	GLY	3.3
1	A	476	GLN	3.3
1	A	507	TYR	3.2
1	B	505	SER	3.2
1	B	525	PHE	3.2
1	B	308	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	613	LEU	3.1
1	B	304	GLU	3.0
1	B	509	TYR	3.0
1	A	321	MET	3.0
1	B	504	ASP	2.9
1	A	509	TYR	2.9
1	A	474	GLN	2.9
1	A	289	GLY	2.9
1	B	474	GLN	2.9
1	A	307	LYS	2.8
1	A	298	GLU	2.8
1	B	414	PRO	2.7
1	A	311	SER	2.7
1	A	508	LEU	2.7
1	B	503	GLN	2.7
1	A	308	ASN	2.6
1	A	332	HIS	2.6
1	B	298	GLU	2.6
1	A	477	GLY	2.6
1	B	301	SER	2.6
1	B	475	MET	2.6
1	A	602	GLU	2.6
1	B	316	LEU	2.6
1	B	317	ASN	2.6
1	A	466	GLN	2.5
1	B	310	ASP	2.5
1	B	473	GLU	2.4
1	B	452	ASN	2.4
1	B	314	ARG	2.4
1	B	598	VAL	2.3
1	B	307	LYS	2.3
1	B	479	LYS	2.3
1	B	356	THR	2.3
1	B	309	ILE	2.3
1	B	599	LYS	2.3
1	B	400	ARG	2.3
1	B	311	SER	2.2
1	B	290	GLU	2.2
1	B	584	GLU	2.2
1	A	317	ASN	2.2
1	A	559	GLN	2.2
1	B	448	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	504	ASP	2.1
1	B	303	GLU	2.1
1	B	293	VAL	2.1
1	A	416	LEU	2.0
1	B	305	TRP	2.0
1	B	477	GLY	2.0
1	A	360	LYS	2.0
1	A	286	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.