



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 02:13 AM GMT

PDB ID : 1D3Q
Title : CRYSTAL STRUCTURE OF HUMAN ALPHA THROMBIN IN COMPLEX
WITH BENZO[B]THIOPHENE INHIBITOR 2
Authors : Chirgadze, N.Y.
Deposited on : 1999-09-30
Resolution : 2.90 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

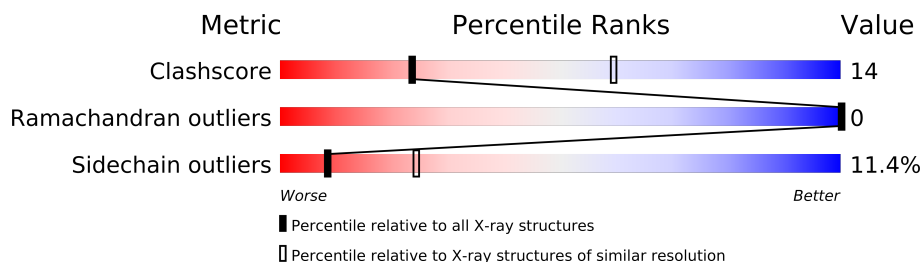
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.90 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))
Clashscore	79885	1326 (2.90-2.90)
Ramachandran outliers	78287	1290 (2.90-2.90)
Sidechain outliers	78261	1292 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	36	
2	B	259	
3	H	12	

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 2418 atoms, of which 0 are hydrogen and 0 are deuterium.

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 ALPHA-THROMBIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	28	Total	C	N	O	S	0	0	0
			231	145	37	48	1			

- Molecule 2 is a protein called ALPHA-THROMBIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	250	Total	C	N	O	S	3	0	0
			2025	1292	358	361	14			

- Molecule 3 is a protein called HIRUGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	12	Total	C	N	O	S	0	0	0
			108	66	13	28	1			

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).

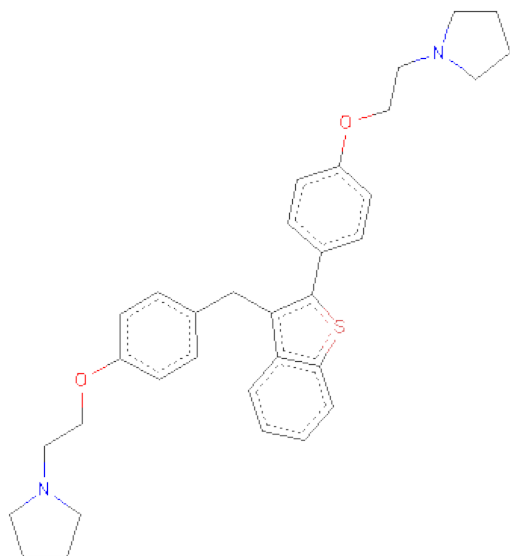


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Na	0	0
			2	2		

- Molecule 6 is 3-[4-(2-PYRROLIDIN-1-YL-ETHOXY)-BENZYL]-2-4-(2-PYRROLIDIN-1-YL-ETHOXY)-PHENYL]-BENZO[B]THIOPHENE (three-letter code: BT2) (formula: C₃₃H₃₈N₂O₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	S	0	0
			38	33	2	2	1		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

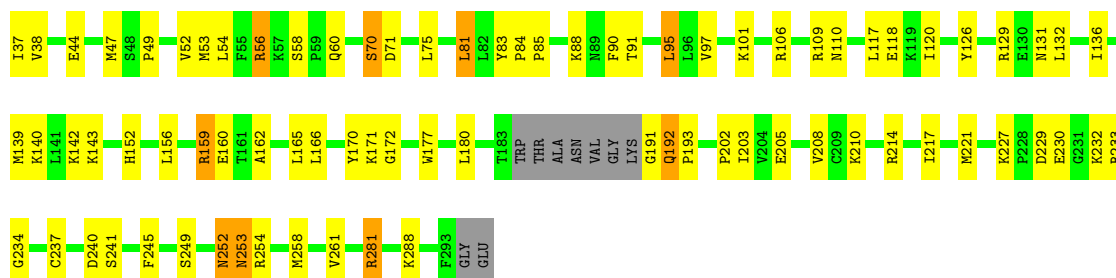
• Molecule 1: ALPHA-THROMBIN

Chain A: 



• Molecule 2: ALPHA-THROMBIN

Chain B: 



• Molecule 3: HIRUGEN

Chain H: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	71.15Å 71.73Å 73.10Å 90.00° 100.56° 90.00°	Depositor
Resolution (Å)	20.00 – 2.90	Depositor
% Data completeness (in resolution range)	85.4 (20.00-2.90)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 98.0	Depositor
R, R_{free}	0.167 , 0.228	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2418	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, BT2, NAG, TYS

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	0.45	0/233	0.67	0/310
2	B	0.45	0/2077	0.68	0/2806
3	H	0.78	0/92	0.74	0/120
All	All	0.47	0/2402	0.68	0/3236

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	231	0	228	11	0
2	B	2025	0	2000	58	0
3	H	108	0	85	1	0
4	B	14	0	13	0	0
5	B	2	0	0	0	0
6	B	38	0	38	2	0
All	All	2418	0	2364	68	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including

hydrogens) of the entry. The overall clashscore for this entry is 14.

All (68) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:13:PRO:HA	1:A:17:LYS:HG3	1.57	0.87
2:B:95:LEU:HD22	2:B:95:LEU:N	2.02	0.74
2:B:95:LEU:HD22	2:B:95:LEU:H	1.54	0.73
2:B:180:LEU:HD21	2:B:193:PRO:HB3	1.70	0.73
1:A:25:GLU:O	1:A:29:LEU:HD23	1.94	0.68
1:A:26:ARG:HH21	1:A:26:ARG:HG2	1.58	0.67
2:B:281:ARG:HH11	2:B:281:ARG:HG3	1.60	0.67
3:H:311:GLN:HG2	3:H:311:GLN:OXT	1.93	0.66
2:B:249:SER:HB3	2:B:252:ASN:ND2	2.11	0.65
2:B:88:LYS:HD2	2:B:90:PHE:HE2	1.63	0.64
2:B:88:LYS:HD2	2:B:90:PHE:CE2	2.34	0.62
2:B:81:LEU:HD23	2:B:120:ILE:HG21	1.81	0.62
2:B:249:SER:HB3	2:B:252:ASN:HD21	1.64	0.61
2:B:95:LEU:HD23	2:B:117:LEU:HD12	1.84	0.58
2:B:131:ASN:HA	6:B:400:BT2:H172	1.85	0.58
2:B:81:LEU:HD23	2:B:120:ILE:CG2	2.34	0.57
2:B:95:LEU:N	2:B:95:LEU:CD2	2.67	0.56
2:B:252:ASN:ND2	2:B:254:ARG:H	2.02	0.56
2:B:56:ARG:O	2:B:60:GLN:HA	2.06	0.55
2:B:53:MET:HB2	2:B:177:TRP:CZ3	2.42	0.54
2:B:109:ARG:O	2:B:110:ASN:HB2	2.08	0.54
2:B:171:LYS:HA	2:B:202:PRO:HA	1.90	0.53
2:B:160:GLU:H	2:B:160:GLU:CD	2.13	0.52
2:B:191:GLY:C	2:B:192:GLN:HE21	2.13	0.52
2:B:131:ASN:HA	6:B:400:BT2:H231	1.91	0.52
2:B:118:GLU:OE2	2:B:140:LYS:HD3	2.11	0.50
2:B:252:ASN:HD22	2:B:252:ASN:C	2.15	0.50
2:B:180:LEU:HD21	2:B:193:PRO:CB	2.39	0.50
1:A:11:LEU:O	2:B:152:HIS:CD2	2.66	0.49
2:B:166:LEU:HD23	2:B:203:ILE:HD13	1.95	0.49
2:B:205:GLU:HB2	2:B:208:VAL:HG23	1.95	0.48
2:B:237:CYS:O	2:B:240:ASP:HB2	2.13	0.48
2:B:227:LYS:HB2	2:B:230:GLU:HG3	1.96	0.47
2:B:159:ARG:HB3	2:B:159:ARG:HH11	1.78	0.47
2:B:252:ASN:C	2:B:252:ASN:ND2	2.68	0.47
2:B:131:ASN:OD1	2:B:131:ASN:N	2.49	0.46
2:B:75:LEU:HD11	2:B:136:ILE:HD11	1.97	0.45
2:B:37:ILE:N	2:B:240:ASP:OD1	2.48	0.45
2:B:252:ASN:HD22	2:B:253:ASN:N	2.14	0.45
1:A:26:ARG:NH2	1:A:26:ARG:HG2	2.30	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:205:GLU:CD	2:B:205:GLU:H	2.20	0.45
2:B:126:TYR:CE1	2:B:132:LEU:HD22	2.52	0.45
2:B:172:GLY:HA3	2:B:245:PHE:CZ	2.51	0.45
2:B:95:LEU:H	2:B:95:LEU:CD2	2.26	0.44
2:B:241:SER:HA	2:B:261:VAL:HB	1.99	0.44
1:A:13:PRO:CA	1:A:17:LYS:HG3	2.38	0.43
2:B:70:SER:OG	2:B:71:ASP:N	2.51	0.43
1:A:6:GLU:HB2	1:A:9:CYS:HB3	1.98	0.43
1:A:12:ARG:HG2	2:B:49:PRO:HG3	1.99	0.43
2:B:101:LYS:HB3	2:B:101:LYS:HE3	1.76	0.43
2:B:81:LEU:N	2:B:81:LEU:CD1	2.81	0.43
2:B:281:ARG:NH1	2:B:281:ARG:HG3	2.32	0.43
2:B:83:TYR:CZ	2:B:85:PRO:HG2	2.53	0.43
1:A:26:ARG:CG	1:A:26:ARG:HH21	2.26	0.43
2:B:38:VAL:O	2:B:234:GLY:HA2	2.18	0.42
2:B:117:LEU:HD22	2:B:139:MET:HB3	2.02	0.42
2:B:281:ARG:HH11	2:B:281:ARG:CG	2.28	0.42
2:B:166:LEU:HA	2:B:166:LEU:HD23	1.87	0.42
2:B:165:LEU:O	2:B:170:TYR:HD2	2.03	0.42
2:B:281:ARG:NH1	2:B:281:ARG:CG	2.83	0.41
2:B:162:ALA:HA	2:B:258:MET:CE	2.51	0.41
1:A:26:ARG:CG	1:A:26:ARG:NH2	2.82	0.41
2:B:95:LEU:CD2	2:B:117:LEU:HD12	2.50	0.41
2:B:84:PRO:O	2:B:85:PRO:C	2.59	0.41
2:B:52:VAL:HG12	2:B:53:MET:N	2.36	0.41
1:A:22:ASP:OD2	1:A:22:ASP:C	2.59	0.41
2:B:44:GLU:O	2:B:47:MET:HB2	2.20	0.41
2:B:129:ARG:HH11	2:B:129:ARG:HD2	1.75	0.40

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	26/36 (72%)	25 (96%)	1 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	246/259 (95%)	227 (92%)	19 (8%)	0	100	100
3	H	9/12 (75%)	9 (100%)	0	0	100	100
All	All	281/307 (92%)	261 (93%)	20 (7%)	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	26/31 (84%)	25 (96%)	1 (4%)	44	83
2	B	219/225 (97%)	194 (89%)	25 (11%)	8	24
3	H	10/10 (100%)	7 (70%)	3 (30%)	0	1
All	All	255/266 (96%)	226 (89%)	29 (11%)	8	24

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ARG
2	B	54	LEU
2	B	56	ARG
2	B	58	SER
2	B	70	SER
2	B	81	LEU
2	B	91	THR
2	B	95	LEU
2	B	97	VAL
2	B	106	ARG
2	B	142	LYS
2	B	143	LYS
2	B	156	LEU
2	B	159	ARG
2	B	192	GLN
2	B	210	LYS
2	B	214	ARG

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Mol	Chain	Res	Type
2	B	217	ILE
2	B	221	MET
2	B	229	ASP
2	B	232	LYS
2	B	233	ARG
2	B	252	ASN
2	B	253	ASN
2	B	281	ARG
2	B	288	LYS
3	H	304	GLU
3	H	308	GLU
3	H	311	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	192	GLN
2	B	252	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	TYS	H	309	3	16,16,17	6.61	14 (87%)	20,22,24	1.78	6 (30%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical

component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	TYS	H	309	3	-	0/9/11/13	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	309	TYS	O-C	18.40	1.24	1.11
3	H	309	TYS	CE2-CD2	6.94	1.52	1.38
3	H	309	TYS	OH-S	-6.92	1.51	1.63
3	H	309	TYS	CE1-CD1	6.56	1.51	1.38
3	H	309	TYS	CE2-CZ	6.35	1.52	1.38
3	H	309	TYS	CE1-CZ	6.25	1.51	1.38
3	H	309	TYS	CD2-CG	6.24	1.53	1.38
3	H	309	TYS	O2-S	6.06	1.67	1.45
3	H	309	TYS	CD1-CG	5.01	1.50	1.38
3	H	309	TYS	O1-S	3.72	1.59	1.45
3	H	309	TYS	OH-CZ	-2.69	1.39	1.42
3	H	309	TYS	O3-S	2.61	1.64	1.50
3	H	309	TYS	CA-C	2.49	1.52	1.48
3	H	309	TYS	CB-CG	-2.22	1.45	1.51

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	309	TYS	CB-CG-CD1	-3.19	114.65	120.90
3	H	309	TYS	C-CA-N	3.18	117.01	113.83
3	H	309	TYS	CZ-OH-S	-3.16	114.24	118.98
3	H	309	TYS	CG-CB-CA	-2.55	108.40	114.42
3	H	309	TYS	OH-CZ-CE2	2.16	122.28	118.70
3	H	309	TYS	CD2-CG-CD1	2.07	121.65	118.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
6	BT2	B	400	-	43,43,43	4.70	24 (55%)	58,58,58	3.48	19 (32%)
4	NAG	B	500	2	12,14,15	0.96	0	15,19,21	1.61	5 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	BT2	B	400	-	-	0/20/34/34	0/4/6/6
4	NAG	B	500	2	-	0/6/23/26	0/1/1/1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	400	BT2	C32-C15	8.23	1.54	1.38
6	B	400	BT2	C6-C5	8.16	1.55	1.36
6	B	400	BT2	C14-C12	8.04	1.54	1.38
6	B	400	BT2	C1-C2	7.84	1.54	1.36
6	B	400	BT2	C2-C3	7.69	1.48	1.40
6	B	400	BT2	C18-C16	7.60	1.54	1.38
6	B	400	BT2	C32-C19	7.10	1.53	1.38
6	B	400	BT2	C3-S1	-7.08	1.67	1.74
6	B	400	BT2	C20-C18	7.05	1.52	1.38
6	B	400	BT2	C14-C16	6.70	1.52	1.38
6	B	400	BT2	C15-C13	6.48	1.53	1.38
6	B	400	BT2	C21-C11	6.48	1.51	1.38
6	B	400	BT2	C21-C19	6.32	1.52	1.38
6	B	400	BT2	C12-C10	5.80	1.51	1.39
6	B	400	BT2	C4-C3	5.60	1.48	1.41
6	B	400	BT2	C11-C13	5.53	1.51	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	400	BT2	C6-C1	5.22	1.53	1.37
6	B	400	BT2	C10-C8	5.19	1.53	1.48
6	B	400	BT2	C7-C9	4.30	1.58	1.52
6	B	400	BT2	C9-C4	3.79	1.46	1.41
6	B	400	BT2	C20-C10	3.45	1.46	1.39
6	B	400	BT2	C22-N3	3.44	1.54	1.47
6	B	400	BT2	C26-N2	3.08	1.53	1.47
6	B	400	BT2	C28-N3	2.79	1.53	1.47

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	400	BT2	C8-S1-C3	15.21	98.29	91.39
6	B	400	BT2	C27-N2-C26	10.82	115.48	104.20
6	B	400	BT2	C28-N3-C22	9.18	113.78	104.20
6	B	400	BT2	C33-N2-C27	5.41	126.39	113.38
6	B	400	BT2	C4-C3-S1	-5.15	109.26	111.69
6	B	400	BT2	C12-C10-C8	4.98	124.77	120.94
6	B	400	BT2	C20-C10-C8	-4.90	117.17	120.94
6	B	400	BT2	C10-C8-S1	4.72	127.17	116.51
6	B	400	BT2	O2-C24-C34	3.92	116.47	107.50
6	B	400	BT2	C9-C4-C3	3.64	111.24	107.11
6	B	400	BT2	O3-C25-C33	3.43	115.34	107.50
4	B	500	NAG	C4-C3-C2	-3.15	103.61	111.32
6	B	400	BT2	C2-C3-C4	-2.83	121.81	123.28
4	B	500	NAG	O3-C3-C4	-2.63	104.46	110.35
6	B	400	BT2	C6-C1-C2	-2.59	116.68	120.47
4	B	500	NAG	C3-C4-C5	-2.38	105.95	110.20
4	B	500	NAG	O4-C4-C5	2.34	115.46	109.28
6	B	400	BT2	C9-C7-C13	2.33	118.77	113.31
4	B	500	NAG	O5-C5-C4	-2.24	107.81	110.65
6	B	400	BT2	C30-C22-N3	-2.18	101.26	103.99
6	B	400	BT2	C11-C21-C19	2.15	122.64	119.76
6	B	400	BT2	C21-C19-C32	-2.10	116.61	120.21
6	B	400	BT2	C9-C8-S1	-2.02	107.71	111.79
6	B	400	BT2	C1-C2-C3	2.02	121.59	118.86

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.