



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 09:38 AM BST

PDB ID : 4UE7  
Title : Thrombin in complex with 1-amidinopiperidine  
Authors : Ruehmann, E.; Heine, A.; Klebe, G.  
Deposited on : 2014-12-16  
Resolution : 1.13 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.13.1

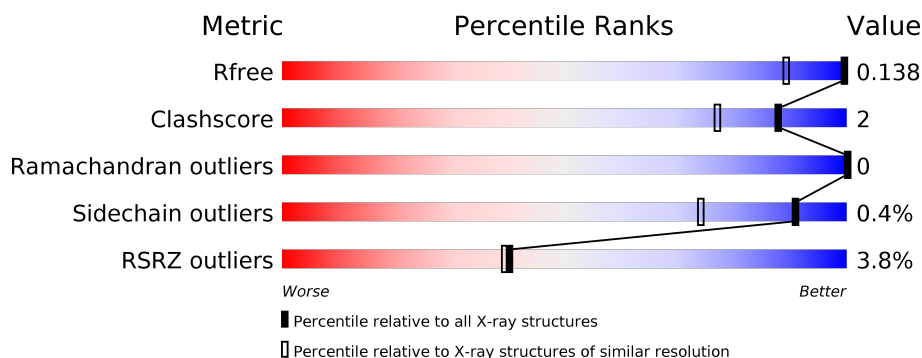
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.13 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))
$R_{free}$	130704	1168 (1.14-1.10)
Clashscore	141614	1205 (1.14-1.10)
Ramachandran outliers	138981	1168 (1.14-1.10)
Sidechain outliers	138945	1165 (1.14-1.10)
RSRZ outliers	127900	1146 (1.14-1.10)

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  $\geq 3$ , 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	H	258	<div> <div>2%</div> <div> <div></div> <div>93%</div> <div>5%</div> </div> </div>
2	I	12	<div> <div>25%</div> <div> <div></div> <div>92%</div> <div>8%</div> </div> </div>
3	L	28	<div> <div>7%</div> <div> <div></div> <div>100%</div> </div> </div>

## 2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 5116 atoms, of which 2302 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 THROMBIN HEAVY CHAIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	H	251	Total	C	H	N	O	S	0	15	0
			4166	1348	2073	363	367	15			

- Molecule 2 is a protein called HIRUDIN VARIANT-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	12	Total	C	N	O	S	0	0	1
			89	57	11	20	1			

- Molecule 3 is a protein called THROMBIN LIGHT CHAIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	L	28	Total	C	H	N	O	S	0	2	0
			465	150	229	36	49	1			

- Molecule 4 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	1	Total	I	0	0
			1	1		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	H	1	Total	C	O	0	0
			6	3	3		
5	H	1	Total	C	O	0	0
			6	3	3		
5	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	H	2	Total	Na	0	0
			2	2		

- Molecule 7 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	H	1	Total	O	P	0	0
			5	4	1		

- Molecule 8 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula:  $C_2H_6OS$ ).



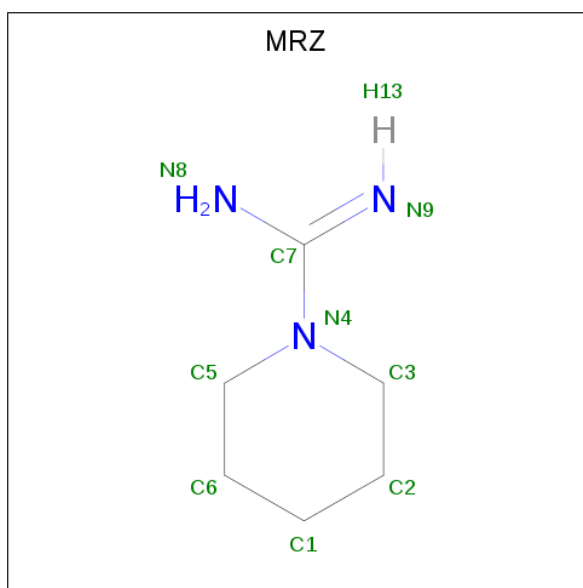
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	H	1	Total	C	O	S	0	0
			4	2	1	1		
8	H	1	Total	C	O	S	0	0
			4	2	1	1		
8	H	1	Total	C	O	S	0	1
			8	4	2	2		

- Molecule 9 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	H	1	Total	C	O	0	0
			4	2	2		

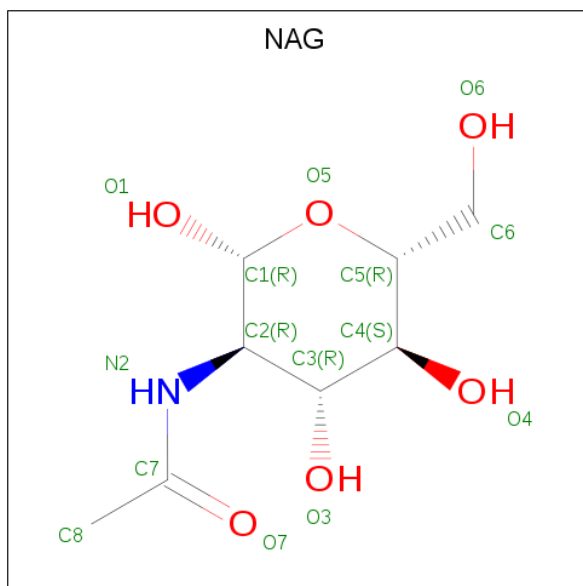
- Molecule 10 is piperidine-1-carboximidamide (three-letter code: MRZ) (formula:  $C_6H_{13}N_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	H	1	Total	C	N	0	0
			9	6	3		

- Molecule 11 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG)

(formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	H	1	Total	C	N	O	0	0
			14	8	1	5		

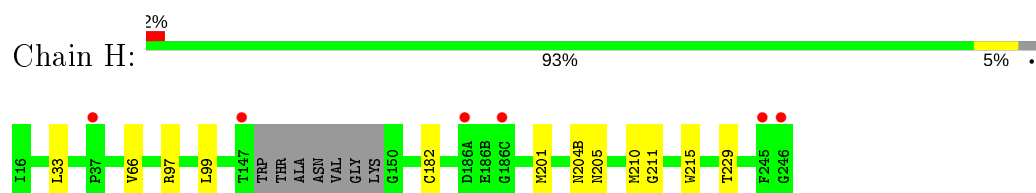
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	H	289	Total	O	0	9
			294	294		
12	I	4	Total	O	0	0
			4	4		
12	L	29	Total	O	0	0
			29	29		

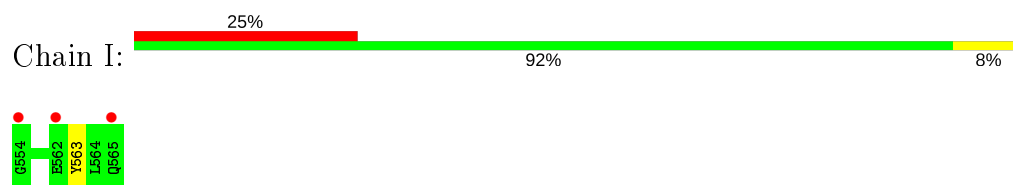
### 3 Residue-property plots

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.

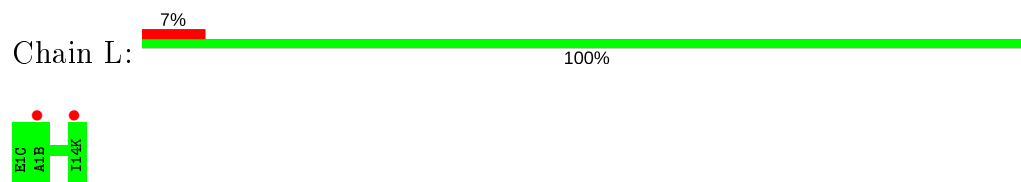
- Molecule 1: THROMBIN HEAVY CHAIN



- Molecule 2: HIRUDIN VARIANT-2



- Molecule 3: THROMBIN LIGHT CHAIN





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.26Å 71.48Å 72.44Å 90.00° 100.30° 90.00°	Depositor
Resolution (Å)	19.56 – 1.13 19.56 – 1.13	Depositor EDS
% Data completeness (in resolution range)	99.5 (19.56-1.13) 99.5 (19.56-1.13)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.44 (at 1.13Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.123 , 0.139 0.124 , 0.138	Depositor DCC
$R_{free}$ test set	6600 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	10.7	Xtriage
Anisotropy	0.315	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.44 , 64.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	5116	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.00% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, NAG, MRZ, NA, PO4, EDO, DMS, IOD, 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	H	0.49	0/2178	0.68	0/2944
2	I	0.31	0/73	0.49	0/96
3	L	0.53	0/244	0.76	0/324
All	All	0.49	0/2495	0.68	0/3364

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 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	H	2093	2073	2076	10	0
2	I	89	0	63	0	0
3	L	236	229	232	0	0
4	H	1	0	0	0	0
5	H	18	0	24	0	0
6	H	2	0	0	0	0
7	H	5	0	0	0	0
8	H	16	0	24	0	0
9	H	4	0	6	0	0
10	H	9	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
11	H	14	0	13	0	0
12	H	294	0	0	3	0
12	I	4	0	0	0	0
12	L	29	0	0	0	0
All	All	2814	2302	2450	10	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 2.

All (10) 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:H:97[B]:ARG:NH2	12:H:2134:HOH:O	1.97	0.94
1:H:97[B]:ARG:CZ	12:H:2134:HOH:O	2.29	0.74
1:H:97[B]:ARG:NE	12:H:2134:HOH:O	2.24	0.71
1:H:33[A]:LEU:CD2	1:H:66:VAL:HG12	2.27	0.64
1:H:201:MET:SD	1:H:210[A]:MET:HG3	2.45	0.56
1:H:33[A]:LEU:HD23	1:H:66:VAL:HG12	1.96	0.46
1:H:204(B):ASN:O	1:H:205[A]:ASN:HB3	2.16	0.45
1:H:211:GLY:HA2	1:H:229:THR:O	2.17	0.44
1:H:204(B):ASN:O	1:H:205[A]:ASN:CB	2.67	0.42
1:H:99:LEU:HD12	1:H:215:TRP:HB3	2.03	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	H	262/258 (102%)	255 (97%)	7 (3%)	0	100	100
2	I	9/12 (75%)	9 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	L	28/28 (100%)	27 (96%)	1 (4%)	0	100	100
All	All	299/298 (100%)	291 (97%)	8 (3%)	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	H	222/224 (99%)	221 (100%)	1 (0%)	88	66
2	I	5/10 (50%)	5 (100%)	0	100	100
3	L	25/26 (96%)	25 (100%)	0	100	100
All	All	252/260 (97%)	251 (100%)	1 (0%)	91	74

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	182	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules 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
2	TYS	I	563	2	15,16,17	1.59	2 (13%)	18,22,24	1.09	1 (5%)

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
2	TYS	I	563	2	-	0/10/11/13	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	563	TYS	OH-S	-4.13	1.52	1.58
2	I	563	TYS	CE1-CD1	2.67	1.43	1.38

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	563	TYS	OH-S-O1	-2.33	100.90	107.71

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 3 are monoatomic - leaving 11 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	$\# Z  > 2$
10	MRZ	H	1258	-	9,9,9	0.71	0	10,11,11	1.19	1 (10%)
11	NAG	H	1259	1	14,14,15	1.95	4 (28%)	17,19,21	1.36	3 (17%)
7	PO4	H	1253	-	4,4,4	0.81	0	6,6,6	0.94	0
8	DMS	H	1255	-	3,3,3	0.56	0	3,3,3	0.18	0
8	DMS	H	1257[A]	-	3,3,3	0.64	0	3,3,3	0.74	0
8	DMS	H	1257[B]	-	3,3,3	0.64	0	3,3,3	0.43	0
5	GOL	H	1248	-	5,5,5	0.42	0	5,5,5	0.35	0
5	GOL	H	1250	-	5,5,5	0.39	0	5,5,5	0.34	0
8	DMS	H	1254	-	3,3,3	0.54	0	3,3,3	0.38	0
9	EDO	H	1256	-	3,3,3	0.38	0	2,2,2	0.42	0
5	GOL	H	1249	-	5,5,5	0.37	0	5,5,5	0.58	0

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
11	NAG	H	1259	1	-	0/6/23/26	0/1/1/1
10	MRZ	H	1258	-	-	0/4/12/12	0/1/1/1
5	GOL	H	1248	-	-	2/4/4/4	-
5	GOL	H	1250	-	-	0/4/4/4	-
9	EDO	H	1256	-	-	0/1/1/1	-
5	GOL	H	1249	-	-	3/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	H	1259	NAG	C7-N2	4.10	1.48	1.34
11	H	1259	NAG	C4-C3	-3.34	1.43	1.52
11	H	1259	NAG	O5-C5	3.04	1.49	1.43
11	H	1259	NAG	C3-C2	-2.50	1.47	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
10	H	1258	MRZ	C2-C3-N4	-2.45	105.82	110.66
11	H	1259	NAG	C8-C7-N2	2.33	120.04	116.10
11	H	1259	NAG	C1-O5-C5	2.30	115.31	112.19
11	H	1259	NAG	O5-C1-C2	2.11	114.61	111.29

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	H	1248	GOL	C1-C2-C3-O3
5	H	1249	GOL	O1-C1-C2-C3
5	H	1249	GOL	O1-C1-C2-O2
5	H	1249	GOL	O2-C2-C3-O3
5	H	1248	GOL	O2-C2-C3-O3

There are no ring outliers.

No monomer is involved in short contacts.

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	H	251/258 (97%)	-0.23	6 (2%) 59 56	7, 13, 28, 45	0
2	I	11/12 (91%)	1.49	3 (27%) 0 1	24, 33, 42, 46	0
3	L	28/28 (100%)	-0.14	2 (7%) 16 16	8, 12, 26, 34	0
All	All	290/298 (97%)	-0.16	11 (3%) 40 39	7, 13, 31, 46	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	565	GLN	5.1
3	L	1(B)	ALA	4.7
1	H	245	PHE	4.7
2	I	554	GLY	4.6
1	H	246	GLY	4.6
3	L	14(K)	ILE	4.3
1	H	147	THR	3.7
1	H	186(C)	GLY	3.5
1	H	186(A)	ASP	3.3
1	H	37	PRO	2.3
2	I	562	GLU	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	TYS	I	563	16/17	0.98	0.07	22,27,34,34	0



### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
11	NAG	H	1259	14/15	0.69	0.27	35,39,42,43	0
5	GOL	H	1249	6/6	0.82	0.21	27,30,33,36	6
5	GOL	H	1250	6/6	0.83	0.18	19,19,20,20	6
5	GOL	H	1248	6/6	0.91	0.11	18,23,24,25	6
9	EDO	H	1256	4/4	0.93	0.12	12,14,14,14	4
8	DMS	H	1257[B]	4/4	0.98	0.10	43,43,43,44	4
10	MRZ	H	1258	9/9	0.98	0.08	10,12,15,15	0
7	PO4	H	1253	5/5	0.98	0.07	12,12,13,14	5
8	DMS	H	1255	4/4	0.98	0.08	19,19,20,21	4
8	DMS	H	1257[A]	4/4	0.98	0.10	12,14,15,18	4
8	DMS	H	1254	4/4	0.99	0.06	10,12,13,13	0
6	NA	H	1252	1/1	1.00	0.03	13,13,13,13	0
4	IOD	H	1247	1/1	1.00	0.03	13,13,13,13	1
6	NA	H	1251	1/1	1.00	0.02	10,10,10,10	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.