



wwPDB X-ray Structure Validation Summary Report

Feb 27, 2014 – 01:20 AM GMT

PDB ID : 2B8O
Title : Crystal Structure of Glu-Gly-Arg-ChloromethylKetone-Factor VIIa/Soluble
Tissue Factor Complex
Authors : Bajaj, S.P.; Schmidt, A.E.; Padmanabhan, K.
Deposited on : 2005-10-08
Resolution : 2.80 Å(reported)

This is a wwPDB validation summary 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>

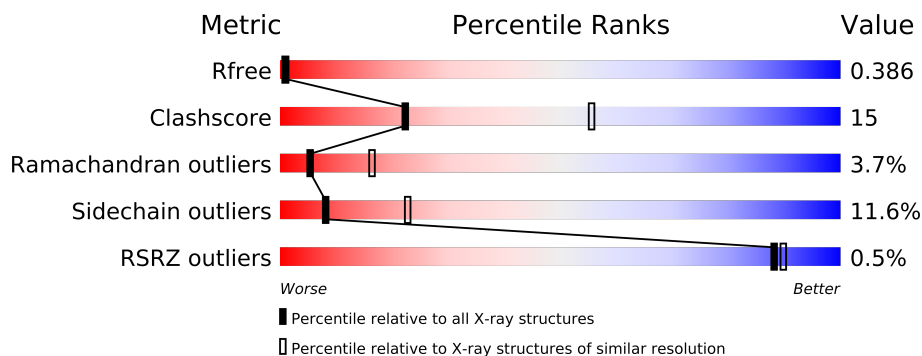
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
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 2.80 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}	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

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.

Mol	Chain	Length	Quality of chain
1	L	142	
2	H	254	
3	T	205	

2 Entry composition i

There are 12 unique types of molecules in this entry. The entry contains 4959 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 Coagulation factor VII light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	142	Total	C	N	O	S	0	0	0
			1135	683	189	248	15			

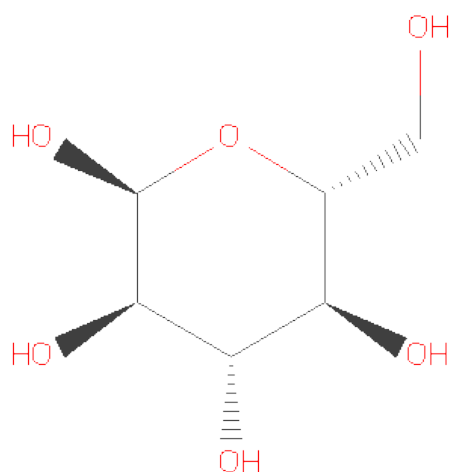
- Molecule 2 is a protein called Coagulation factor VII heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	254	Total	C	N	O	S	0	0	0
			1974	1253	351	357	13			

- Molecule 3 is a protein called Tissue factor.

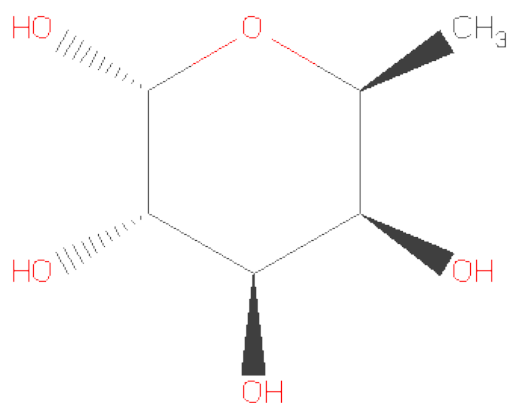
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	T	191	Total	C	N	O	S	0	0	0
			1551	987	250	309	5			

- Molecule 4 is SUGAR (ALPHA-D-GLUCOSE) (three-letter code: GLC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	L	1	Total	C	O	0	0
			11	6	5		

- Molecule 5 is SUGAR (ALPHA-L-FUCOSE) (three-letter code: FUC) (formula: C₆H₁₂O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	L	1	Total	C	O	0	0
			10	6	4		

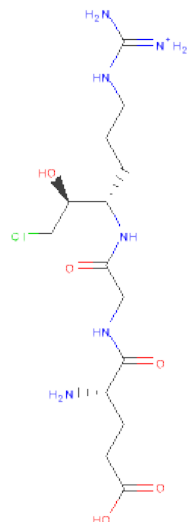
- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	H	1	Total	Ca	0	0
			1	1		
6	L	5	Total	Ca	0	0
			5	5		

- Molecule 7 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	L	2	Total	Mg	0	0
			2	2		

- Molecule 8 is L-ALPHA-GLUTAMYL-N-{(1S)-4-[AMINO(IMINIO)METHYL]AMINO}-1-[(1S)-2-CHLORO-1-HYDROXYETHYL]BUTYL}GLYCINAMIDE (three-letter code: 0GJ) (formula: C₁₄H₂₈ClN₆O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	H	1	Total	C	N	O	0	0
			25	14	6	5		

- Molecule 9 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	H	1	Total	Na	0	0
			1	1		

- Molecule 10 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	H	2	Total	Zn	0	0
			2	2		

- Molecule 11 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	H	2	Total	Cl	0	0
			2	2		
11	T	1	Total	Cl	0	0
			1	1		

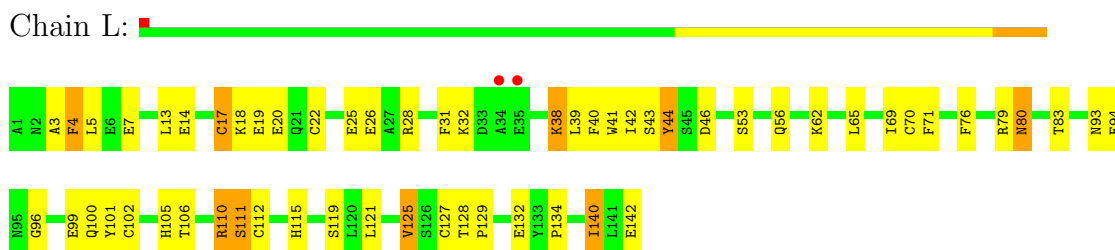
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	L	66	Total 66	O 66	0	0
12	H	91	Total 91	O 91	0	0
12	T	82	Total 82	O 82	0	0

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.

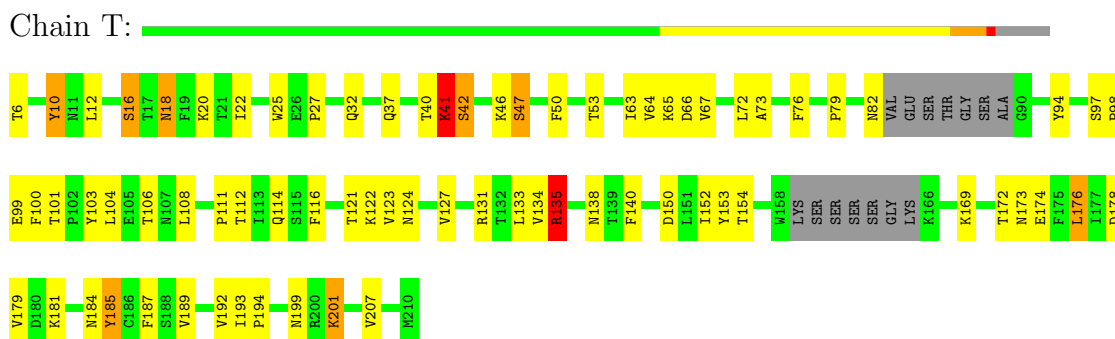
- Molecule 1: Coagulation factor VII light chain



- Molecule 2: Coagulation factor VII heavy chain



- Molecule 3: Tissue factor



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.28Å 81.11Å 126.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.80 40.72 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.80) 54.3 (40.72-2.20)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.73 (at 2.20Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.225 , 0.285 0.234 , 0.386	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	24.8	Xtriage
Anisotropy	0.944	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 37.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 20351 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4959	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, CL, NA, CA, GLC, OGJ, MG, FUC, CGU

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	L	0.64	0/1029	0.93	1/1374 (0.1%)
2	H	0.72	0/2024	0.99	1/2755 (0.0%)
3	T	0.72	0/1585	0.99	3/2156 (0.1%)
All	All	0.70	0/4638	0.97	5/6285 (0.1%)

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 outliers	#Planarity outliers
1	L	0	1
3	T	0	1
All	All	0	2

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	T	135	ARG	NE-CZ-NH1	6.15	123.37	120.30
3	T	20	LYS	N-CA-C	-5.21	96.93	111.00
1	L	17	CYS	CA-CB-SG	5.15	123.26	114.00
2	H	209	LEU	CA-CB-CG	5.07	126.96	115.30
3	T	66	ASP	N-CA-C	-5.04	97.39	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	44	TYR	Sidechain
3	T	103	TYR	Sidechain

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	L	1135	0	987	38	0
2	H	1974	0	1950	60	0
3	T	1551	0	1501	46	0
4	L	11	0	10	0	0
5	L	10	0	10	3	0
6	H	1	0	0	0	0
6	L	5	0	0	0	0
7	L	2	0	0	0	0
8	H	25	0	25	8	0
9	H	1	0	0	0	0
10	H	2	0	0	0	0
11	H	2	0	0	4	0
11	T	1	0	0	1	0
12	H	91	0	0	1	0
12	L	66	0	0	0	0
12	T	82	0	0	3	0
All	All	4959	0	4483	137	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 15.

The worst 5 of 137 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
8:H:1:OGJ:HB3	11:H:514:CL:CL	2.03	0.94
2:H:196:GLY:HA2	2:H:212:ILE:HG23	1.57	0.84
2:H:81:GLN:OE1	2:H:112:VAL:HG12	1.77	0.83
3:T:10:TYR:HD1	3:T:10:TYR:H	1.31	0.79
2:H:44:GLY:HA2	2:H:196:GLY:O	1.86	0.75

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	L	130/142 (92%)	108 (83%)	18 (14%)	4 (3%)	7	21
2	H	252/254 (99%)	220 (87%)	23 (9%)	9 (4%)	5	17
3	T	185/205 (90%)	160 (86%)	17 (9%)	8 (4%)	4	13
All	All	567/601 (94%)	488 (86%)	58 (10%)	21 (4%)	5	16

5 of 21 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	48	ASN
2	H	195	SER
2	H	248	PRO
3	T	41	LYS
3	T	42	SER

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	L	114/114 (100%)	99 (87%)	15 (13%)	6	16
2	H	216/216 (100%)	192 (89%)	24 (11%)	9	25
3	T	178/189 (94%)	158 (89%)	20 (11%)	9	25
All	All	508/519 (98%)	449 (88%)	59 (12%)	8	23

5 of 59 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	H	139	SER
2	H	192	LYS
3	T	176	LEU
2	H	145	LEU
2	H	159	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
2	H	63	ASN
2	H	109	HIS
2	H	143	GLN
2	H	40	GLN
2	H	117	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

10 non-standard protein/DNA/RNA residues are 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	$\# Z > 2$
1	CGU	L	14	1,7	11,11,12	5.43	1 (9%)	12,14,16	3.11	2 (16%)
1	CGU	L	16	1,7,6	11,11,12	5.66	3 (27%)	12,14,16	2.77	2 (16%)
1	CGU	L	19	1,7	11,11,12	5.31	4 (36%)	12,14,16	2.56	5 (41%)
1	CGU	L	20	1	11,11,12	5.31	5 (45%)	12,14,16	2.81	5 (41%)
1	CGU	L	25	1,6	11,11,12	5.20	1 (9%)	12,14,16	1.86	2 (16%)
1	CGU	L	26	1,7	11,11,12	5.83	2 (18%)	12,14,16	2.05	1 (8%)
1	CGU	L	29	1,6	11,11,12	5.85	4 (36%)	12,14,16	2.10	3 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	CGU	L	35	1	11,11,12	5.47	2 (18%)	12,14,16	1.88	2 (16%)
1	CGU	L	6	1,6	11,11,12	5.86	4 (36%)	12,14,16	1.62	3 (25%)
1	CGU	L	7	1,6	11,11,12	6.53	3 (27%)	12,14,16	1.74	4 (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
1	CGU	L	14	1,7	-	0/12/14/16	0/0/0/0
1	CGU	L	16	1,7,6	-	0/12/14/16	0/0/0/0
1	CGU	L	19	1,7	-	0/12/14/16	0/0/0/0
1	CGU	L	20	1	-	0/12/14/16	0/0/0/0
1	CGU	L	25	1,6	-	0/12/14/16	0/0/0/0
1	CGU	L	26	1,7	-	0/12/14/16	0/0/0/0
1	CGU	L	29	1,6	-	0/12/14/16	0/0/0/0
1	CGU	L	35	1	-	0/12/14/16	0/0/0/0
1	CGU	L	6	1,6	-	0/12/14/16	0/0/0/0
1	CGU	L	7	1,6	-	0/12/14/16	0/0/0/0

The worst 5 of 29 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	7	CGU	O-C	20.66	1.25	1.11
1	L	26	CGU	O-C	18.87	1.24	1.11
1	L	29	CGU	O-C	18.64	1.24	1.11
1	L	6	CGU	O-C	18.38	1.24	1.11
1	L	16	CGU	O-C	18.22	1.23	1.11

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	14	CGU	C-CA-N	-9.23	104.61	113.83
1	L	16	CGU	C-CA-N	8.00	121.82	113.83
1	L	19	CGU	C-CA-N	-6.53	107.30	113.83
1	L	26	CGU	C-CA-N	-6.23	107.61	113.83
1	L	29	CGU	C-CA-N	-5.48	108.35	113.83

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 17 ligands modelled in this entry, 14 are monoatomic - leaving 3 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	$\# Z > 2$
8	0GJ	H	1	-	23,24,25	1.81	8 (34%)	28,30,31	1.41	4 (14%)
4	GLC	L	501	1	10,11,12	1.33	1 (10%)	11,15,17	0.78	0
5	FUC	L	502	1	9,10,11	1.09	1 (11%)	10,14,16	0.71	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
8	0GJ	H	1	-	-	0/29/29/31	0/0/0/0
4	GLC	L	501	1	1/1/4/5	0/2/19/22	0/1/1/1
5	FUC	L	502	1	1/1/4/5	0/0/17/20	0/1/1/1

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	H	1	0GJ	C2-CA2	3.29	1.58	1.53
8	H	1	0GJ	CA1-C1	-2.98	1.42	1.52
8	H	1	0GJ	CZ-NE	2.62	1.38	1.33
8	H	1	0GJ	CZ-NH2	2.58	1.38	1.32
8	H	1	0GJ	CB1-CA2	2.46	1.56	1.53

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	H	1	0GJ	CB-CA-N	3.17	117.88	110.14
8	H	1	0GJ	CB1-CA2-C2	2.49	116.74	113.02
8	H	1	0GJ	C2-CA2-N2	-2.49	105.45	110.29
8	H	1	0GJ	O-C-N1	-2.04	118.90	123.05

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	L	502	FUC	C1
4	L	501	GLC	C1

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 ⓘ

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	L	142/142 (100%)	-0.67	2 (1%) 72 72	3, 21, 38, 46	0
2	H	254/254 (100%)	-0.74	0 100 100	2, 12, 28, 47	0
3	T	191/205 (93%)	-0.89	0 100 100	3, 17, 34, 40	0
All	All	587/601 (97%)	-0.77	2 (0%) 88 93	2, 15, 34, 47	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	35	CGU	2.4
1	L	34	ALA	2.3

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	CGU	L	25	12/13	0.19	-	20,33,37,42	0
1	CGU	L	6	12/13	0.22	-	27,32,39,40	0
1	CGU	L	35	12/13	0.23	-	40,46,50,51	0
1	CGU	L	14	12/13	0.14	-	26,41,43,43	0
1	CGU	L	26	12/13	0.11	-	22,27,36,37	0
1	CGU	L	19	12/13	0.13	-	29,38,46,47	0
1	CGU	L	16	12/13	0.10	-	9,12,26,28	0
1	CGU	L	20	12/13	0.15	-	30,37,45,45	0
1	CGU	L	7	12/13	0.15	-	21,27,29,30	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	CGU	L	29	12/13	0.12	-	18,22,26,26	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	CA	L	509	1/1	0.06	-	29,29,29,29	0
11	CL	T	516	1/1	0.10	-	23,23,23,23	0
7	MG	L	505	1/1	0.13	-	25,25,25,25	0
11	CL	H	514	1/1	0.31	-	49,49,49,49	0
6	CA	L	503	1/1	0.13	-	50,50,50,50	0
6	CA	L	507	1/1	0.14	-	66,66,66,66	0
10	ZN	H	512	1/1	0.22	-	44,44,44,44	0
7	MG	L	508	1/1	0.06	-	38,38,38,38	0
5	FUC	L	502	10/11	0.20	-	34,36,38,39	0
4	GLC	L	501	11/12	0.36	-	44,48,51,51	0
11	CL	H	515	1/1	0.03	-	27,27,27,27	0
6	CA	L	504	1/1	0.15	-	60,60,60,60	0
8	OGJ	H	1	25/26	0.26	-	20,27,34,34	0
6	CA	L	506	1/1	0.03	-	55,55,55,55	0
6	CA	H	510	1/1	0.12	-	35,35,35,35	0
9	NA	H	511	1/1	0.43	-	49,49,49,49	0
10	ZN	H	513	1/1	0.44	-	57,57,57,57	0

6.5 Other polymers ⓘ

There are no such residues in this entry.