



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 07:04 PM GMT

PDB ID : 1DVA
Title : Crystal Structure of the Complex Between the Peptide Exosite Inhibitor E-76 and Coagulation Factor VIIA
Authors : Eigenbrot, C.; Ultsch, M.H.
Deposited on : 2000-01-20
Resolution : 3.00 Å(reported)

This is a wwPDB X-ray Structure 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/validation/2016/XrayValidationReportHelp>
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:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

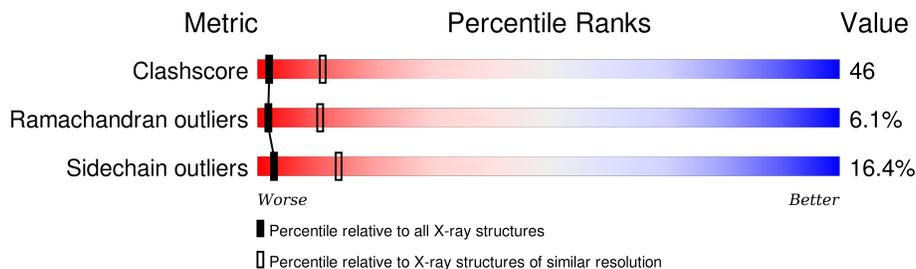
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	H	254	
1	I	254	
2	L	101	
2	M	101	
3	X	20	
3	Y	20	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
9	GLC	M	503	X	-	-	-

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 5917 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 DES-GLA FACTOR VIIA (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	H	254	1974	1253	351	357	13	85	0	0
1	I	254	1974	1253	351	357	13	118	0	0

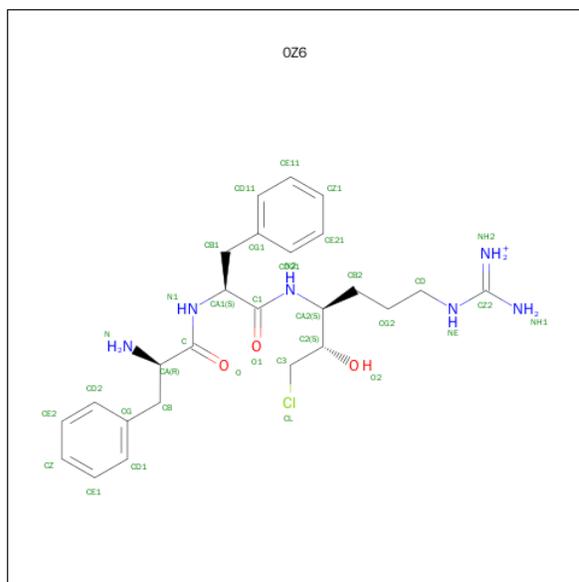
- Molecule 2 is a protein called DES-GLA FACTOR VIIA (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	101	759	456	130	160	13	42	0	0
2	M	101	759	456	130	160	13	116	0	0

- Molecule 3 is a protein called PEPTIDE E-76.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	X	19	151	95	26	28	2	0	0	1
3	Y	19	151	95	26	28	2	0	0	1

- Molecule 4 is D-PHENYLALANYL-N-[(2S,3S)-6- {[AMINO(IMINIO)METHYL]AMINO}-1-CHLORO-2-HYDROXYHEXAN-3-YL]-L-PHENYLALANINAMIDE (three-letter code: 0Z6) (formula: C₂₅H₃₆ClN₆O₃).

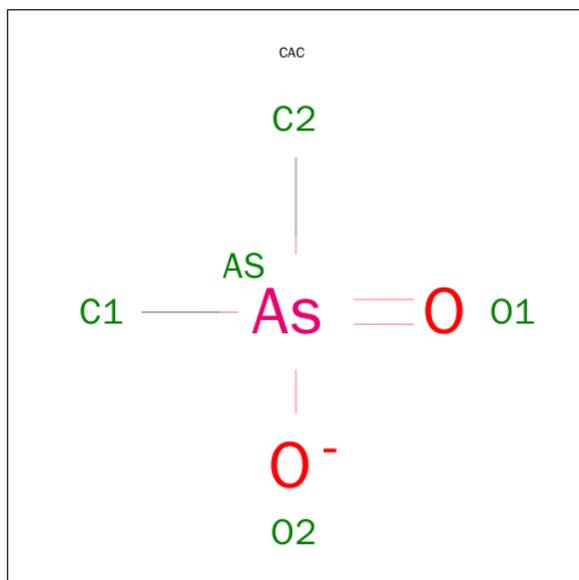


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	H	1	Total	C	N	O	0	0
			34	25	6	3		
4	I	1	Total	C	N	O	0	0
			34	25	6	3		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 6 is CACODYLATE ION (three-letter code: CAC) (formula: C₂H₆AsO₂).

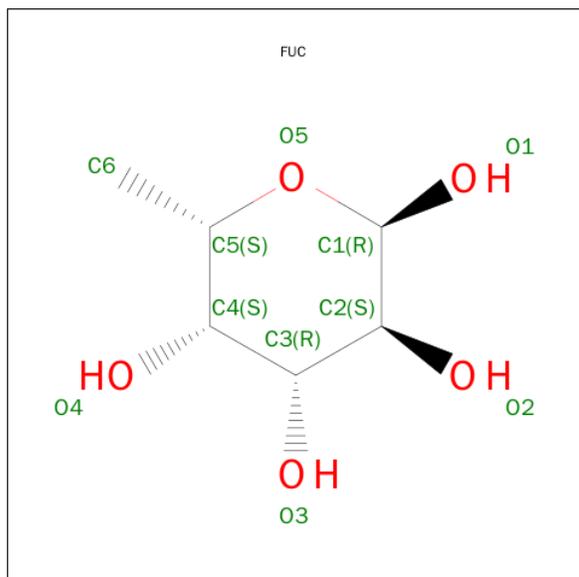


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	As	C	O		
6	H	1	5	1	2	2	4	0
6	I	1	5	1	2	2	4	0

- Molecule 7 is a polymer of unknown type called SUGAR (2-MER).

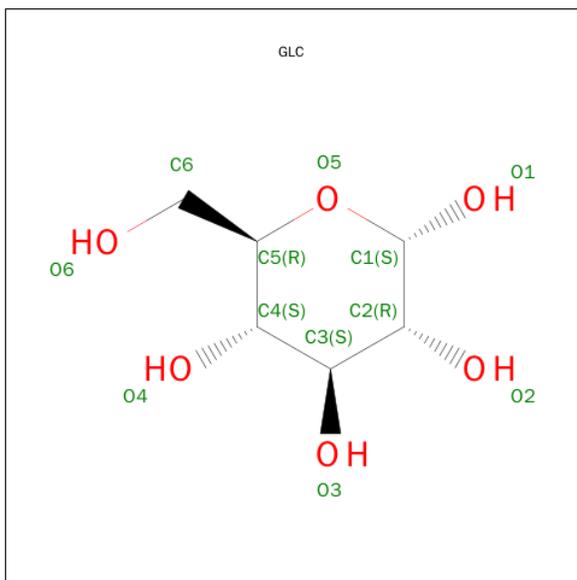
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
7	L	2	22	12	10	0	0

- Molecule 8 is SUGAR (ALPHA-L-FUCOSE) (three-letter code: FUC) (formula: C₆H₁₂O₅).



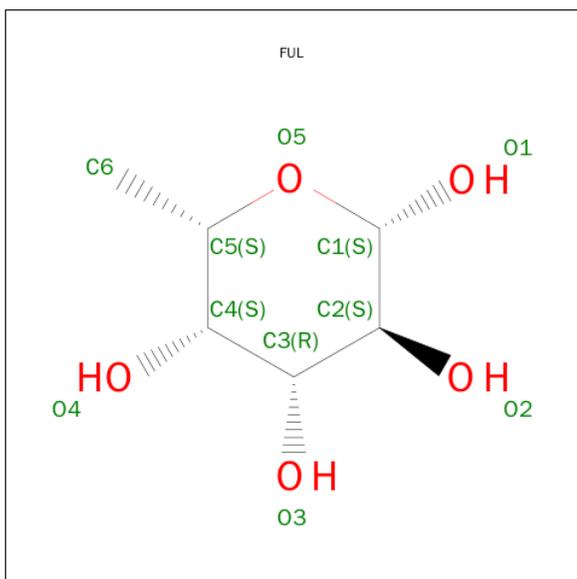
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	L	1	Total	C	O	0	0
			10	6	4		
8	L	1	Total	C	O	0	0
			10	6	4		

- Molecule 9 is SUGAR (ALPHA-D-GLUCOSE) (three-letter code: GLC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	M	1	Total	C	O	0	0
			11	6	5		

- Molecule 10 is SUGAR (BETA-L-FUCOSE) (three-letter code: FUL) (formula: C₆H₁₂O₅).



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

- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	H	3	Total	O	0	0
			3	3		
11	I	1	Total	O	0	0
			1	1		



- Molecule 2: DES-GLA FACTOR VIIA (LIGHT CHAIN)



- Molecule 3: PEPTIDE E-76



- Molecule 3: PEPTIDE E-76



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.49Å 55.26Å 111.73Å 90.00° 99.48° 90.00°	Depositor
Resolution (Å)	50.00 – 3.00	Depositor
% Data completeness (in resolution range)	97.5 (50.00-3.00)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.225 , 0.295	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	5917	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BGC, ACE, CA, GLC, GAL, FUC, 0Z6, CAC, FUL

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	0.63	0/2024	0.87	2/2755 (0.1%)
1	I	0.60	0/2024	0.85	1/2755 (0.0%)
2	L	0.53	0/773	0.75	0/1043
2	M	0.55	0/773	0.74	0/1043
3	X	0.76	0/153	1.05	1/209 (0.5%)
3	Y	0.73	0/153	1.04	1/209 (0.5%)
All	All	0.61	0/5900	0.84	5/8014 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	X	1	ALA	N-CA-C	-6.96	92.20	111.00
1	H	170(C)	ARG	N-CA-C	-6.89	92.39	111.00
3	Y	1	ALA	N-CA-C	-6.55	93.31	111.00
1	H	199	HIS	N-CA-C	-5.44	96.32	111.00
1	I	199	HIS	N-CA-C	-5.42	96.36	111.00

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	H	1974	0	1949	195	0
1	I	1974	0	1949	197	0
2	L	759	0	679	60	0
2	M	759	0	679	48	0
3	X	151	0	132	18	0
3	Y	151	0	132	18	0
4	H	34	0	32	4	0
4	I	34	0	32	3	0
5	H	1	0	0	0	0
5	I	1	0	0	0	0
5	L	1	0	0	0	0
5	M	1	0	0	0	0
6	H	5	0	0	0	0
6	I	5	0	0	0	0
7	L	22	0	19	3	0
8	L	20	0	20	2	0
9	M	11	0	10	0	0
10	M	10	0	10	0	0
11	H	3	0	0	0	0
11	I	1	0	0	1	0
All	All	5917	0	5643	496	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 46.

The worst 5 of 496 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:247:ARG:HG3	1:H:248:PRO:CD	1.82	1.09
1:H:232:SER:HA	1:H:235:ILE:HD11	1.36	1.03
1:H:247:ARG:HG3	1:H:248:PRO:HD2	1.04	1.02
1:I:232:SER:HA	1:I:235:ILE:HD11	1.40	1.00
1:H:247:ARG:CG	1:H:248:PRO:HD2	1.95	0.95

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 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 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	252/254 (99%)	202 (80%)	34 (14%)	16 (6%)	2	9
1	I	252/254 (99%)	200 (79%)	40 (16%)	12 (5%)	3	17
2	L	99/101 (98%)	83 (84%)	8 (8%)	8 (8%)	1	5
2	M	99/101 (98%)	81 (82%)	11 (11%)	7 (7%)	1	7
3	X	17/20 (85%)	15 (88%)	1 (6%)	1 (6%)	2	11
3	Y	17/20 (85%)	14 (82%)	2 (12%)	1 (6%)	2	11
All	All	736/750 (98%)	595 (81%)	96 (13%)	45 (6%)	2	11

5 of 45 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	49	THR
1	H	78	GLY
1	H	146	ASP
1	H	170(A)	GLN
2	L	123	ASP

5.3.2 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 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	216/216 (100%)	178 (82%)	38 (18%)	2	12
1	I	216/216 (100%)	179 (83%)	37 (17%)	2	12
2	L	88/88 (100%)	77 (88%)	11 (12%)	6	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	M	88/88 (100%)	77 (88%)	11 (12%)	6	24
3	X	16/16 (100%)	12 (75%)	4 (25%)	1	3
3	Y	16/16 (100%)	12 (75%)	4 (25%)	1	3
All	All	640/640 (100%)	535 (84%)	105 (16%)	3	14

5 of 105 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	L	137	LYS
1	I	64	LEU
2	M	120	LEU
2	L	140	ILE
1	I	24	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
1	I	30	GLN
1	I	60(D)	ASN
1	I	175	ASN
3	X	14	GLN
1	I	217	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](#)

2 carbohydrates 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
7	BGC	L	501	2,7	11,11,12	0.83	0	14,15,17	1.54	1 (7%)
7	GAL	L	502	7	11,11,12	0.89	0	14,15,17	0.93	1 (7%)

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
7	BGC	L	501	2,7	-	0/2/19/22	0/1/1/1
7	GAL	L	502	7	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	L	502	GAL	C1-C2-C3	-2.54	106.54	109.54
7	L	501	BGC	C1-C2-C3	4.63	115.02	109.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	L	501	BGC	3	0
7	L	502	GAL	2	0

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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
4	OZ6	H	1	1	31,35,36	1.69	1 (3%)	39,45,46	0.85	3 (7%)
6	CAC	H	310	-	0,4,4	0.00	-	0,6,6	0.00	-
4	OZ6	I	1	1	31,35,36	1.75	1 (3%)	39,45,46	0.96	3 (7%)
6	CAC	I	311	-	0,4,4	0.00	-	0,6,6	0.00	-
8	FUC	L	504	2	10,10,11	1.02	1 (10%)	14,14,16	1.27	1 (7%)
8	FUC	L	506	-	10,10,11	0.69	0	14,14,16	0.78	0
9	GLC	M	503	2	11,11,12	0.81	0	14,15,17	0.61	0
10	FUL	M	505	2	10,10,11	0.48	0	14,14,16	0.74	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
4	OZ6	H	1	1	-	0/33/35/37	0/2/2/2
6	CAC	H	310	-	-	0/0/0/0	0/0/0/0
4	OZ6	I	1	1	-	0/33/35/37	0/2/2/2
6	CAC	I	311	-	-	0/0/0/0	0/0/0/0
8	FUC	L	504	2	-	0/0/17/20	0/1/1/1
8	FUC	L	506	-	-	0/0/17/20	0/1/1/1
9	GLC	M	503	2	1/1/4/5	0/2/19/22	0/1/1/1
10	FUL	M	505	2	-	0/0/17/20	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	I	1	OZ6	O2-C2	-9.22	1.22	1.43
4	H	1	OZ6	O2-C2	-9.17	1.22	1.43
8	L	504	FUC	C1-C2	-2.14	1.47	1.52

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	1	OZ6	CA2-N2-C1	-2.78	117.97	123.13
4	H	1	OZ6	CA2-N2-C1	-2.53	118.42	123.13
8	L	504	FUC	O2-C2-C1	-2.44	104.32	109.21
4	H	1	OZ6	C1-CA1-N1	-2.04	105.51	111.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	1	0Z6	C2-CA2-N2	2.46	114.71	110.11

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
9	M	503	GLC	C1

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	1	0Z6	4	0
4	I	1	0Z6	3	0
8	L	504	FUC	2	0

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

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.