



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 04:44 PM BST

PDB ID : 2VDR  
Title : Integrin AlphaIIbBeta3 Headpiece Bound to a chimeric Fibrinogen Gamma chain peptide, LGGAKQRGDV  
Authors : Springer, T.A.; Zhu, J.; Xiao, T.  
Deposited on : 2007-10-10  
Resolution : 2.40 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

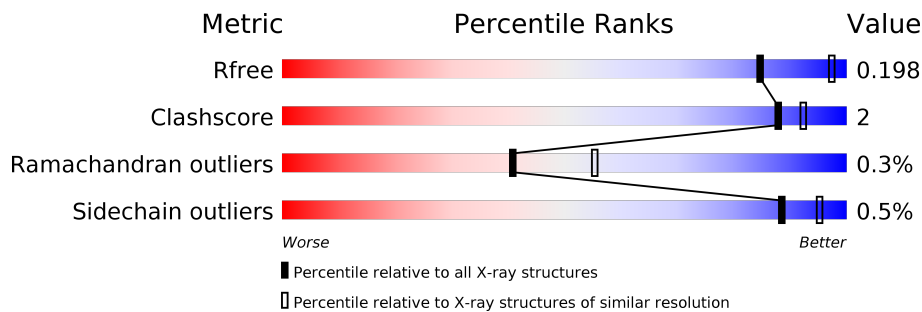
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)

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  $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	452	96% (Green), 4% (Yellow), 0% (Grey), 0% (Red), 0% (Orange)
2	B	461	94% (Green), 5% (Yellow), 1% (Grey), 0% (Red), 0% (Orange)
3	C	10	70% (Green), 30% (Grey)
4	H	221	94% (Green), 5% (Yellow), 1% (Grey), 0% (Red), 0% (Orange)
5	L	214	94% (Green), 6% (Yellow), 0% (Grey), 0% (Red), 0% (Orange)
6	D	5	60% (Green), 40% (Yellow)
7	E	7	43% (Green), 57% (Yellow)

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:

<b>Mol</b>	<b>Type</b>	<b>Chain</b>	<b>Res</b>	<b>Chirality</b>	<b>Geometry</b>	<b>Clashes</b>	<b>Electron density</b>
6	MAN	D	3	X	-	-	-
7	MAN	E	3	X	-	-	-

## 2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 11884 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 INTEGRIN ALPHA-IIB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	452	3539	2250	616	665	8	0	9	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	282	GLY	ALA	conflict	UNP P08514

- Molecule 2 is a protein called INTEGRIN BETA-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	455	3580	2236	607	704	33	0	10	0

- Molecule 3 is a protein called FIBRINOGEN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	7	54	31	12	11	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	408	ARG	ALA	engineered mutation	UNP Q53Y18

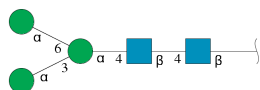
- Molecule 4 is a protein called MONOCLONAL ANTIBODY 10E5 HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	H	219	1678	1064	270	338	6	0	3	0

- Molecule 5 is a protein called MONOCLONAL ANTIBODY 10E5 LIGHT CHAIN.

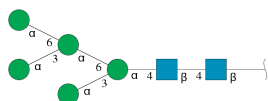
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	L	214	1686	1051	273	351	11	0	9	0

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
6	D	5	61	34	2	25		0	0	0

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
7	E	7	83	46	2	35		0	0	0

- Molecule 8 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
8	A	1	Total C O 6 3 3	0	0
8	B	1	Total C O 6 3 3	0	0

- Molecule 9 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	B	2	Total Ca 2 2	0	0
9	A	4	Total Ca 4 4	0	0

- Molecule 10 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
10	A	1	Total	C	N	O	0	0
			14	8	1	5		
10	A	1	Total	C	N	O	0	0
			14	8	1	5		
10	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 11 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	B	1	Total	Mg	0	0
			1	1		

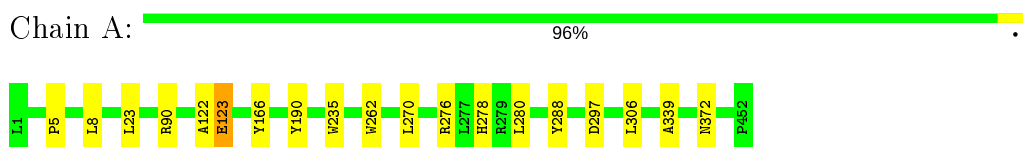
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	412	Total	O	0	0
			412	412		
12	B	256	Total	O	0	0
			256	256		
12	C	12	Total	O	0	0
			12	12		
12	H	214	Total	O	0	0
			214	214		
12	L	248	Total	O	0	0
			248	248		

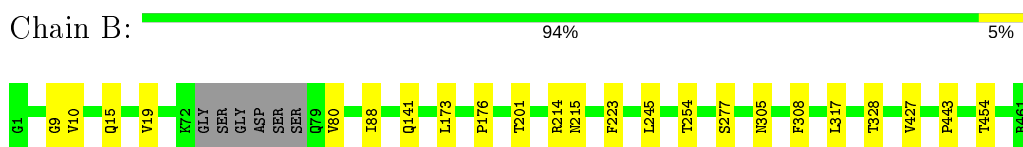
### 3 Residue-property plots [i](#)

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. 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. 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.

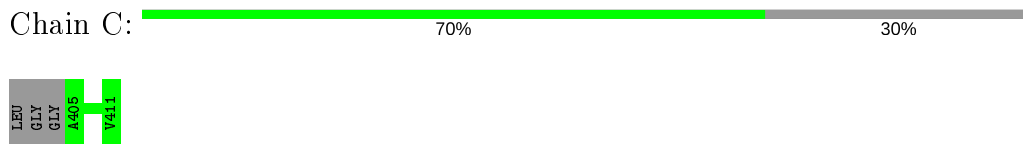
- Molecule 1: INTEGRIN ALPHA-IIB



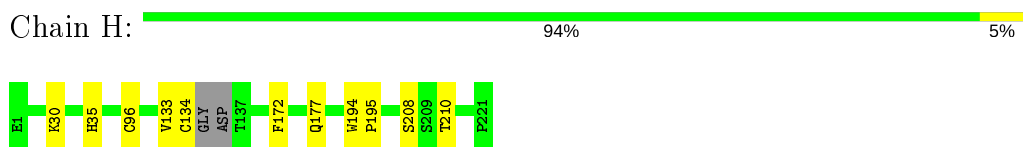
- Molecule 2: INTEGRIN BETA-3



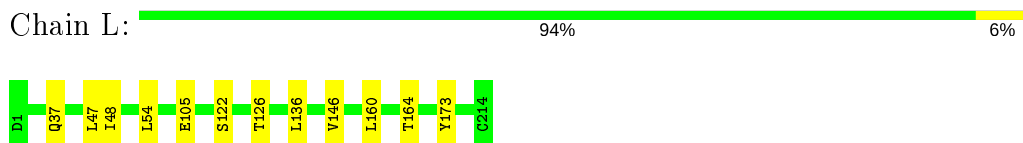
- Molecule 3: FIBRINOGEN



- Molecule 4: MONOCLONAL ANTIBODY 10E5 HEAVY CHAIN



- Molecule 5: MONOCLONAL ANTIBODY 10E5 LIGHT CHAIN



- Molecule 6: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

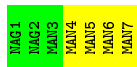






- Molecule 7:  $\alpha$ -D-mannopyranose-(1-3)-[ $\alpha$ -D-mannopyranose-(1-6)] $\alpha$ -D-mannopyranose-(1-6)-[ $\alpha$ -D-mannopyranose-(1-3)] $\alpha$ -D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	148.33Å 148.33Å 176.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.44 – 2.40 43.42 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.5 (43.44-2.40) 96.2 (43.42-2.40)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.02 (at 2.39Å)	Xtrriage
Refinement program	REFMAC 5.3.0037	Depositor
R, $R_{free}$	0.148 , 0.193 0.154 , 0.198	Depositor DCC
$R_{free}$ test set	4276 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.2	Xtrriage
Anisotropy	0.037	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 42.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.019 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	11884	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	55.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, MG, CA, NAG, MAN

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	0.37	0/3656	0.57	0/4980
2	B	0.34	0/3661	0.50	0/4960
3	C	0.41	0/53	0.61	0/67
4	H	0.35	0/1730	0.53	0/2369
5	L	0.36	0/1754	0.56	0/2377
All	All	0.36	0/10854	0.54	0/14753

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 [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	3539	0	3385	12	0
2	B	3580	0	3523	12	0
3	C	54	0	54	0	0
4	H	1678	0	1637	13	0
5	L	1686	0	1614	14	0
6	D	61	0	52	0	0
7	E	83	0	70	0	0
8	A	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	B	6	0	7	0	0
9	A	4	0	0	0	0
9	B	2	0	0	0	0
10	A	28	0	26	0	0
10	B	14	0	13	0	0
11	B	1	0	0	0	0
12	A	412	0	0	1	0
12	B	256	0	0	4	0
12	C	12	0	0	0	0
12	H	214	0	0	2	0
12	L	248	0	0	0	0
All	All	11884	0	10389	43	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 (43) 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)
4:H:177[B]:GLN:OE1	5:L:160[B]:LEU:HD11	1.46	1.15
4:H:177[B]:GLN:OE1	5:L:160[B]:LEU:CD1	2.07	1.01
1:A:270:LEU:HD23	1:A:276[A]:ARG:HA	1.81	0.63
4:H:177[B]:GLN:CD	5:L:160[B]:LEU:CD1	2.68	0.61
1:A:270:LEU:HD23	1:A:276[B]:ARG:HA	1.82	0.60
2:B:201[A]:THR:HG22	12:B:4097:HOH:O	2.02	0.59
4:H:177[B]:GLN:CD	5:L:160[B]:LEU:HD11	2.21	0.59
5:L:48:ILE:HD13	5:L:54:LEU:HD12	1.85	0.58
5:L:136:LEU:HD21	5:L:146:VAL:HG22	1.84	0.58
4:H:177[B]:GLN:OE1	5:L:160[B]:LEU:HD12	1.99	0.57
4:H:177[A]:GLN:NE2	12:H:4197:HOH:O	2.32	0.54
4:H:30:LYS:NZ	12:H:4046:HOH:O	2.41	0.52
1:A:280:LEU:CD1	1:A:306:LEU:HD23	2.41	0.51
5:L:122:SER:O	5:L:126:THR:HG23	2.10	0.51
5:L:48:ILE:CD1	5:L:54:LEU:HD12	2.41	0.50
2:B:173:LEU:HD12	12:B:4068:HOH:O	2.11	0.50
2:B:88[B]:ILE:HD11	2:B:427:VAL:HG22	1.93	0.50
5:L:136:LEU:HD12	5:L:136:LEU:N	2.27	0.48
4:H:133:VAL:HG12	4:H:134:CYS:N	2.29	0.48
1:A:280:LEU:CD1	1:A:306:LEU:CD2	2.93	0.47
1:A:122:ALA:O	1:A:123:GLU:HB2	2.14	0.47
2:B:141:GLN:NE2	12:B:4053:HOH:O	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:308:PHE:CE2	2:B:328:THR:HG21	2.50	0.46
2:B:201[A]:THR:HG23	12:B:4095:HOH:O	2.15	0.46
5:L:37:GLN:HB2	5:L:47:LEU:HD11	1.98	0.46
1:A:278[A]:HIS:CD2	1:A:339:ALA:HB1	2.51	0.46
5:L:105[A]:GLU:HG3	5:L:173:TYR:OH	2.17	0.45
1:A:235:TRP:CZ2	1:A:270:LEU:HD11	2.52	0.45
2:B:223:PHE:CZ	2:B:254:THR:HG21	2.52	0.44
1:A:90[A]:ARG:CZ	12:A:4053:HOH:O	2.64	0.44
4:H:194:TRP:CG	4:H:195:PRO:HA	2.52	0.44
1:A:5:PRO:HA	1:A:8:LEU:HD21	2.00	0.44
4:H:208:SER:O	4:H:210:THR:HG23	2.18	0.43
2:B:443:PRO:HA	2:B:454:THR:HG22	2.01	0.43
1:A:278[A]:HIS:NE2	1:A:339:ALA:HB1	2.34	0.43
2:B:245:LEU:HD23	2:B:305:ASN:HB2	2.00	0.43
4:H:172:PHE:CD1	5:L:164:THR:HG23	2.54	0.42
1:A:297:ASP:O	1:A:372:ASN:HB2	2.20	0.42
1:A:262:TRP:HB3	2:B:317:LEU:HD13	2.01	0.42
2:B:15:GLN:O	2:B:19:VAL:HG23	2.20	0.41
4:H:177[B]:GLN:HB3	5:L:160[B]:LEU:HD11	2.02	0.41
4:H:35:HIS:O	4:H:96:CYS:HA	2.21	0.41
2:B:176:PRO:HB2	2:B:214:ARG:HB3	2.03	0.40

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	A	459/452 (102%)	450 (98%)	8 (2%)	1 (0%)	47 62
2	B	460/461 (100%)	440 (96%)	17 (4%)	3 (1%)	22 32
3	C	5/10 (50%)	4 (80%)	1 (20%)	0	100 100
4	H	218/221 (99%)	213 (98%)	5 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
5	L	223/214 (104%)	218 (98%)	5 (2%)	0	100 100
All	All	1365/1358 (100%)	1325 (97%)	36 (3%)	4 (0%)	41 55

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	GLU
2	B	9	GLY
2	B	10	VAL
2	B	80	VAL

### 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	A	369/360 (102%)	365 (99%)	4 (1%)	73 87
2	B	414/409 (101%)	412 (100%)	2 (0%)	88 95
3	C	5/6 (83%)	5 (100%)	0	100 100
4	H	192/190 (101%)	192 (100%)	0	100 100
5	L	199/188 (106%)	199 (100%)	0	100 100
All	All	1179/1153 (102%)	1173 (100%)	6 (0%)	88 95

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LEU
1	A	166	TYR
1	A	190	TYR
1	A	288	TYR
2	B	215	ASN
2	B	277	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	316	ASN

### 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](#)

12 monosaccharides 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	$\# Z  > 2$
6	NAG	D	1	2,6	14,14,15	0.59	0	17,19,21	0.91	1 (5%)
6	NAG	D	2	6	14,14,15	0.69	0	17,19,21	0.81	0
6	MAN	D	3	6	11,11,12	0.53	0	15,15,17	0.68	0
6	MAN	D	4	6	11,11,12	0.55	0	15,15,17	0.70	0
6	MAN	D	5	6	11,11,12	0.63	0	15,15,17	1.21	2 (13%)
7	NAG	E	1	2,7	14,14,15	0.73	0	17,19,21	1.00	0
7	NAG	E	2	7	14,14,15	0.56	0	17,19,21	0.85	0
7	MAN	E	3	7	11,11,12	0.66	0	15,15,17	0.71	0
7	MAN	E	4	7	11,11,12	0.50	0	15,15,17	1.19	1 (6%)
7	MAN	E	5	7	11,11,12	0.53	0	15,15,17	0.98	1 (6%)
7	MAN	E	6	7	11,11,12	0.59	0	15,15,17	0.78	1 (6%)
7	MAN	E	7	7	11,11,12	0.59	0	15,15,17	0.81	1 (6%)

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	NAG	D	1	2,6	-	0/6/23/26	0/1/1/1
6	NAG	D	2	6	-	0/6/23/26	0/1/1/1
6	MAN	D	3	6	1/1/4/5	0/2/19/22	0/1/1/1
6	MAN	D	4	6	-	2/2/19/22	0/1/1/1
6	MAN	D	5	6	-	2/2/19/22	0/1/1/1
7	NAG	E	1	2,7	-	0/6/23/26	0/1/1/1
7	NAG	E	2	7	-	2/6/23/26	0/1/1/1
7	MAN	E	3	7	1/1/4/5	0/2/19/22	0/1/1/1
7	MAN	E	4	7	-	0/2/19/22	0/1/1/1
7	MAN	E	5	7	-	0/2/19/22	0/1/1/1
7	MAN	E	6	7	-	2/2/19/22	0/1/1/1
7	MAN	E	7	7	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	E	4	MAN	C1-O5-C5	3.59	117.06	112.19
6	D	5	MAN	C1-C2-C3	3.30	113.72	109.67
7	E	5	MAN	C1-O5-C5	2.56	115.66	112.19
7	E	7	MAN	C1-O5-C5	2.16	115.11	112.19
7	E	6	MAN	C1-O5-C5	2.06	114.99	112.19
6	D	1	NAG	O5-C1-C2	-2.04	108.06	111.29
6	D	5	MAN	C1-O5-C5	2.01	114.92	112.19

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
7	E	3	MAN	C1
6	D	3	MAN	C1

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	5	MAN	O5-C5-C6-O6
7	E	6	MAN	O5-C5-C6-O6
7	E	6	MAN	C4-C5-C6-O6
7	E	2	NAG	O5-C5-C6-O6
6	D	5	MAN	C4-C5-C6-O6
6	D	4	MAN	C4-C5-C6-O6
7	E	2	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
6	D	4	MAN	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

## 5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 7 are monoatomic - leaving 5 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	GOL	A	1453	-	5,5,5	0.44	0	5,5,5	0.11	0
8	GOL	B	1462	9	5,5,5	0.40	0	5,5,5	0.31	0
10	NAG	A	3249	1	14,14,15	0.48	0	17,19,21	0.98	1 (5%)
10	NAG	A	3015	1	14,14,15	0.48	0	17,19,21	0.78	0
10	NAG	B	3099	2	14,14,15	0.50	0	17,19,21	0.90	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
8	GOL	A	1453	-	-	1/4/4/4	-
8	GOL	B	1462	9	-	2/4/4/4	-
10	NAG	A	3249	1	-	0/6/23/26	0/1/1/1
10	NAG	A	3015	1	-	0/6/23/26	0/1/1/1
10	NAG	B	3099	2	-	2/6/23/26	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( $^{\circ}$ )	Ideal( $^{\circ}$ )
10	A	3249	NAG	O5-C1-C2	-2.32	107.63	111.29
10	B	3099	NAG	O5-C1-C2	-2.11	107.96	111.29

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	B	1462	GOL	O1-C1-C2-C3
10	B	3099	NAG	O5-C5-C6-O6
10	B	3099	NAG	C4-C5-C6-O6
8	B	1462	GOL	O1-C1-C2-O2
8	A	1453	GOL	O1-C1-C2-C3

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

### 6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.