



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 10:34 PM GMT

PDB ID : 1U8C  
Title : A novel adaptation of the integrin PSI domain revealed from its crystal structure  
Authors : Xiong, J.P.; Stehle, T.; Goodman, S.L.; Arnaout, M.A.  
Deposited on : 2004-08-05  
Resolution : 3.10 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](mailto: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.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (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

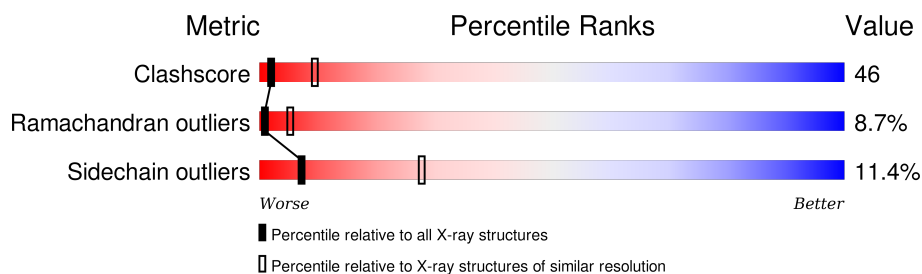
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)

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. 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	A	957	
2	B	692	

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
3	NAG	A	2044	-	-	X	-
3	NAG	A	2045	-	-	X	-

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 12115 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-V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	927	Total	C	N	O	S	0	0	0
			7216	4568	1224	1389	35			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	753	VAL	ILE	SEE REMARK 999	UNP P06756

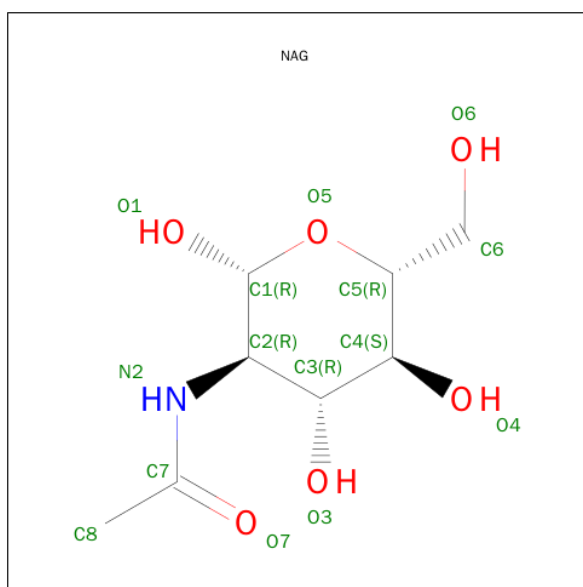
- Molecule 2 is a protein called Integrin beta-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	599	Total	C	N	O	S	0	0	0
			4627	2864	777	930	56			

- Molecule 3 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	2	Total	C	N	O	0	0
			28	16	2	10		
3	A	2	Total	C	N	O	0	0
			28	16	2	10		
3	A	2	Total	C	N	O	0	0
			28	16	2	10		

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

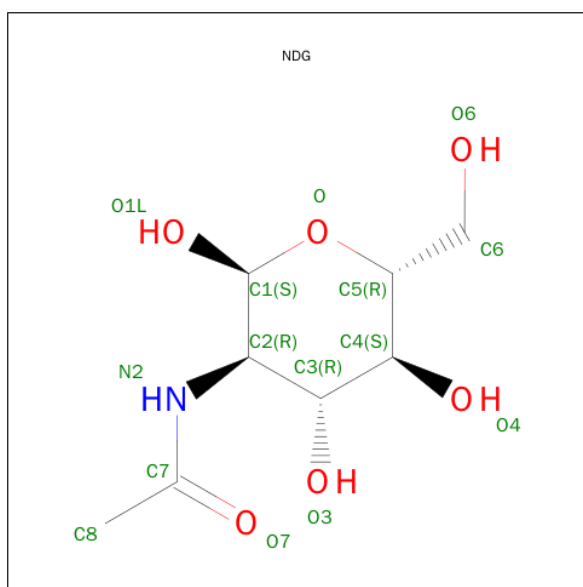


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 6 is SUGAR (2-(ACETYLAMINO)-2-DEOXY-A-D-GLUCOPYRANOSE) (three-letter code: NDG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



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

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	2	Total	C	N	O	0	0
			28	16	2	10		
7	B	2	Total	C	N	O	0	0
			28	16	2	10		
7	B	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 8 is CALCIUM ION (three-letter code: CA) (formula: Ca).

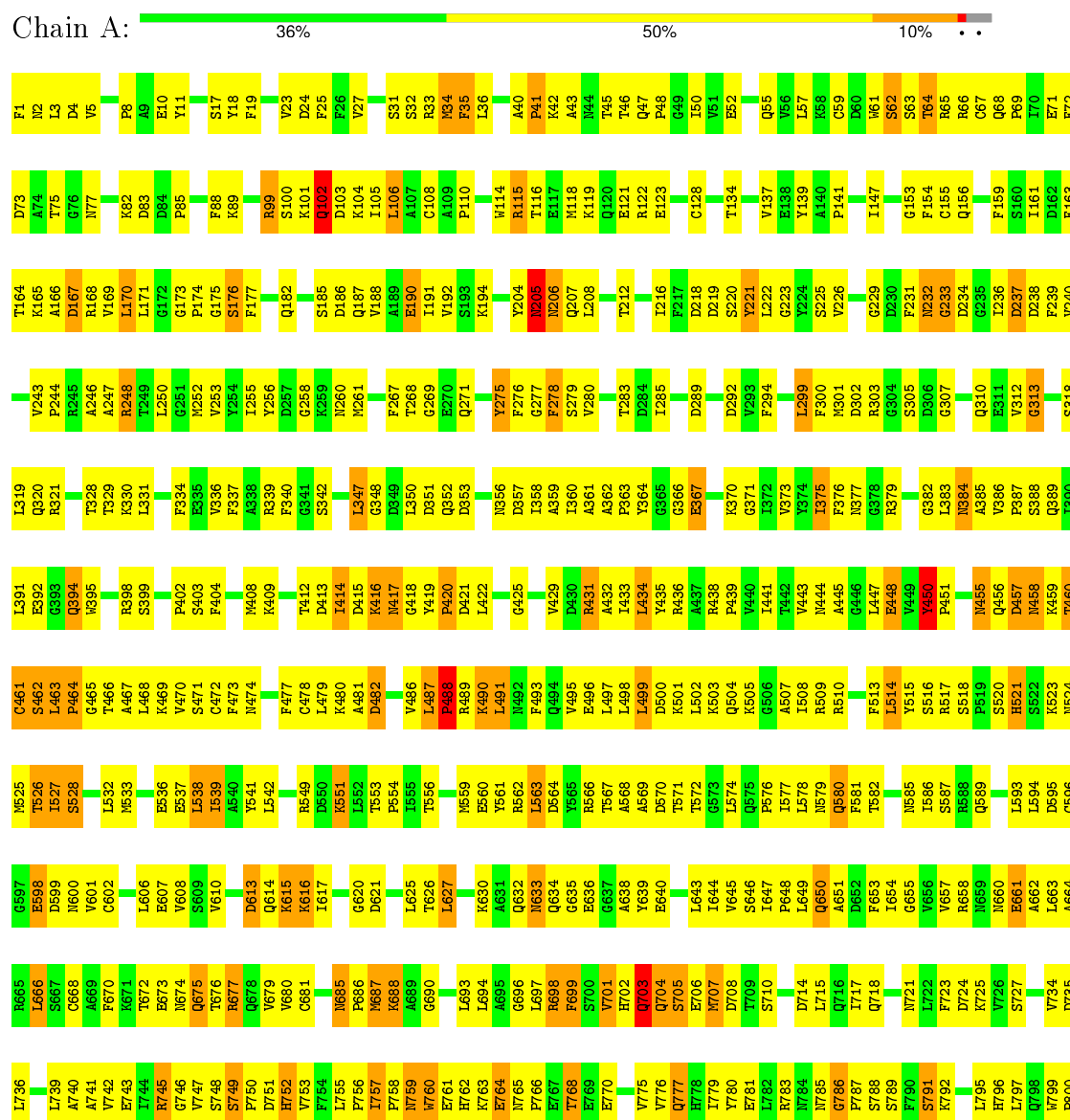
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total	Ca	0	0
			1	1		
8	A	5	Total	Ca	0	0
			5	5		

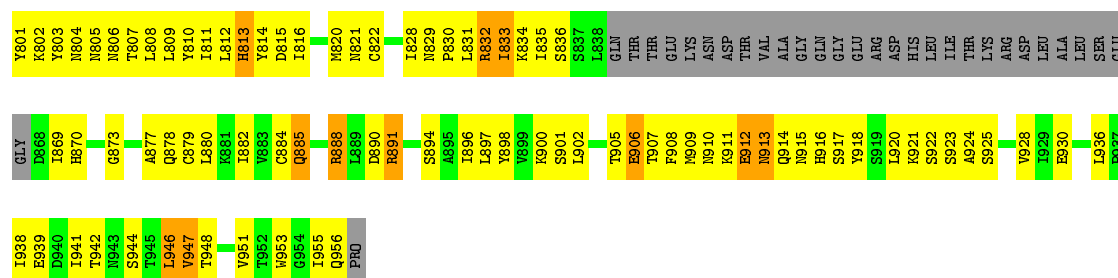
### 3 Residue-property plots

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 ( $\text{RSRZ} > 2$ ). 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.

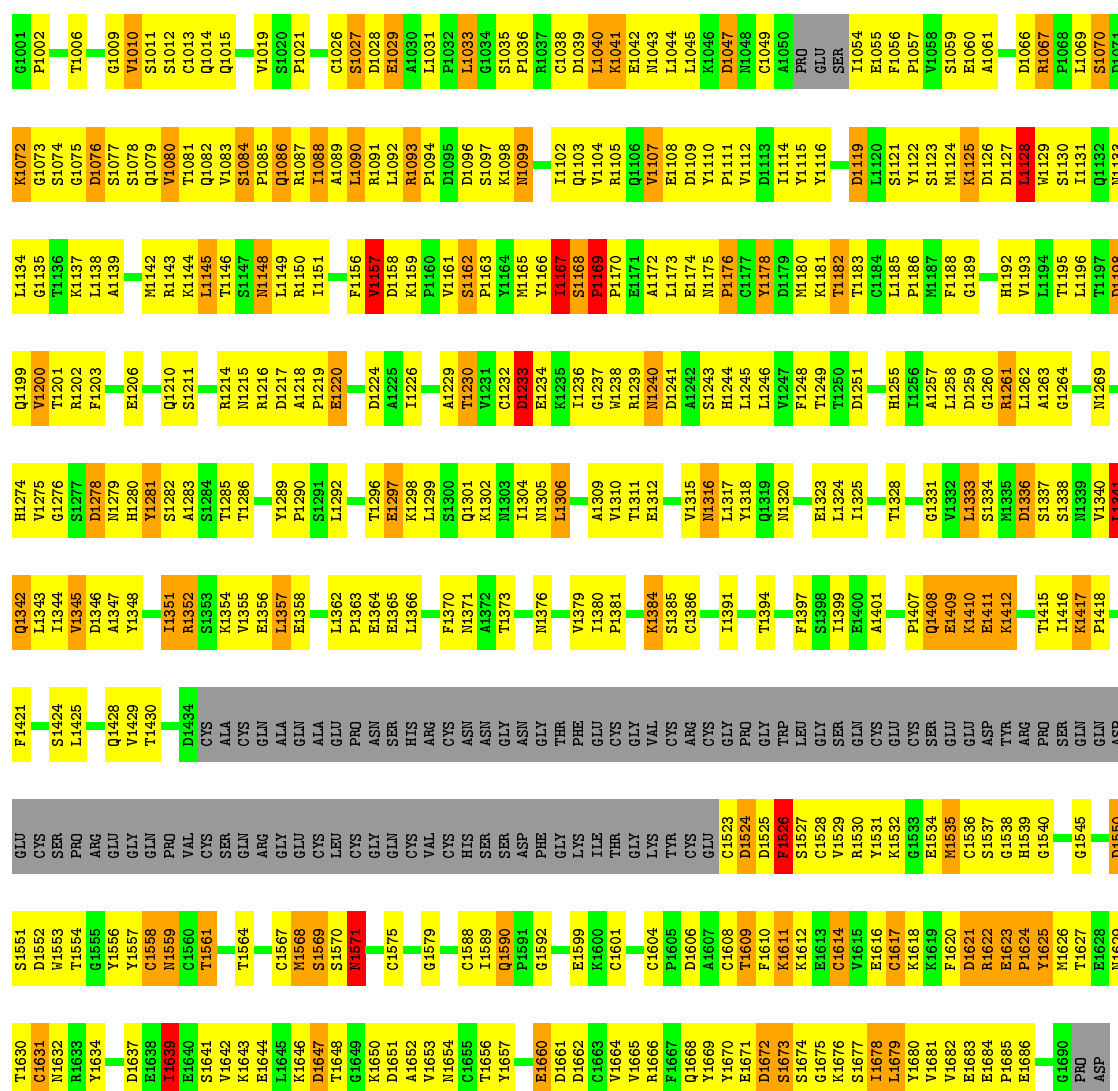
#### • Molecule 1: Integrin alpha-V





• Molecule 2: Integrin beta-3

Chain B: 31% 43% 11% 13%



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 3 2 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	130.00Å 130.00Å 307.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.10	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-3.10)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.293 , 0.367	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	12115	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	55.0	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: CA, NAG, NDG

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	0.46	0/7372	0.71	0/9994
2	B	0.43	0/4708	0.71	1/6367 (0.0%)
All	All	0.45	0/12080	0.71	1/16361 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1169	PRO	C-N-CD	6.01	141.02	128.40

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	7216	0	7033	637	0
2	B	4627	0	4449	450	0
3	A	84	0	75	10	0
4	A	28	0	26	9	0
4	B	14	0	13	2	0
5	A	28	0	25	2	0
6	A	14	0	13	0	0
6	B	14	0	13	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	28	0	25	0	0
7	B	56	0	50	9	0
8	A	5	0	0	0	0
8	B	1	0	0	0	0
All	All	12115	0	11722	1100	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 1100 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)
2:B:1072:LYS:HE2	2:B:1075:GLY:HA3	1.30	1.14
1:A:487:LEU:HD23	1:A:488:PRO:HD3	1.24	1.11
2:B:1041:LYS:HD2	2:B:1045:LEU:CD1	1.80	1.11
2:B:1040:LEU:O	2:B:1042:GLU:N	1.86	1.09
2:B:1611:LYS:HG3	2:B:1669:TYR:HE2	1.17	1.08

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	923/957 (96%)	700 (76%)	161 (17%)	62 (7%)	1	9
2	B	593/692 (86%)	403 (68%)	120 (20%)	70 (12%)	0	2
All	All	1516/1649 (92%)	1103 (73%)	281 (18%)	132 (9%)	1	5

5 of 132 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	SER
1	A	102	GLN
1	A	233	GLY
1	A	414	ILE
1	A	416	LYS

### 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	787/812 (97%)	698 (89%)	89 (11%)	7	28
2	B	537/616 (87%)	475 (88%)	62 (12%)	7	27
All	All	1324/1428 (93%)	1173 (89%)	151 (11%)	7	28

5 of 151 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	745	ARG
1	A	911	LYS
2	B	1611	LYS
1	A	764	GLU
1	A	832	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 52 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	704	GLN
1	A	798	GLN
2	B	1428	GLN
1	A	718	GLN
1	A	777	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

14 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	$\# Z  > 2$
3	NAG	A	2044	1,3	14,14,15	0.53	0	15,19,21	0.95	1 (6%)
3	NAG	A	2045	3	14,14,15	0.59	0	15,19,21	0.86	1 (6%)
3	NAG	A	2266	1,3	14,14,15	0.51	0	15,19,21	0.85	1 (6%)
3	NAG	A	2267	3	14,14,15	0.44	0	15,19,21	0.78	1 (6%)
5	NDG	A	2585	1,5	14,14,15	0.74	0	15,19,21	1.09	2 (13%)
5	NAG	A	2586	5	14,14,15	0.50	0	15,19,21	0.72	1 (6%)
7	NAG	A	2943	1,7	14,14,15	0.63	0	15,19,21	1.04	1 (6%)
7	NDG	A	2944	7	14,14,15	0.63	0	15,19,21	0.78	1 (6%)
3	NAG	A	2950	1,3	14,14,15	0.66	0	15,19,21	1.02	1 (6%)
3	NAG	A	2951	3	14,14,15	0.58	0	15,19,21	0.60	0
7	NAG	B	3559	2,7	14,14,15	0.75	0	15,19,21	1.15	3 (20%)
7	NDG	B	3560	7	14,14,15	0.74	1 (7%)	15,19,21	0.73	1 (6%)
7	NAG	B	3654	2,7	14,14,15	0.76	0	15,19,21	1.31	1 (6%)
7	NDG	B	3655	7	14,14,15	0.65	0	15,19,21	0.87	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
3	NAG	A	2044	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	2045	3	-	0/6/23/26	0/1/1/1
3	NAG	A	2266	1,3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	2267	3	-	0/6/23/26	0/1/1/1
5	NDG	A	2585	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	2586	5	-	0/6/23/26	0/1/1/1
7	NAG	A	2943	1,7	-	0/6/23/26	0/1/1/1
7	NDG	A	2944	7	-	0/6/23/26	0/1/1/1
3	NAG	A	2950	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	2951	3	-	0/6/23/26	0/1/1/1
7	NAG	B	3559	2,7	-	0/6/23/26	0/1/1/1
7	NDG	B	3560	7	-	0/6/23/26	0/1/1/1
7	NAG	B	3654	2,7	-	0/6/23/26	0/1/1/1
7	NDG	B	3655	7	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	3560	NDG	C1-C2	2.14	1.55	1.52

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2950	NAG	C2-N2-C7	-3.15	118.99	123.04
7	A	2943	NAG	C2-N2-C7	-2.83	119.40	123.04
3	A	2266	NAG	C2-N2-C7	-2.79	119.46	123.04
7	B	3655	NDG	C2-N2-C7	-2.57	119.73	123.04
3	A	2267	NAG	C2-N2-C7	-2.44	119.90	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2044	NAG	8	0
3	A	2045	NAG	8	0
5	A	2585	NDG	2	0
5	A	2586	NAG	1	0
3	A	2950	NAG	2	0
3	A	2951	NAG	1	0
7	B	3559	NAG	4	0
7	B	3560	NDG	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	3654	NAG	3	0
7	B	3655	NDG	2	0

## 5.6 Ligand geometry

Of 11 ligands modelled in this entry, 6 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	$\# Z  > 2$
4	NAG	A	2260	1	14,14,15	0.61	0	15,19,21	0.83	0
4	NAG	A	2458	1	14,14,15	0.66	0	15,19,21	0.69	0
6	NDG	A	2821	1	14,14,15	0.67	0	15,19,21	0.77	0
6	NDG	B	3320	2	14,14,15	0.82	0	15,19,21	0.92	1 (6%)
4	NAG	B	3371	2	14,14,15	0.72	0	15,19,21	0.88	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
4	NAG	A	2260	1	-	0/6/23/26	0/1/1/1
4	NAG	A	2458	1	-	0/6/23/26	0/1/1/1
6	NDG	A	2821	1	-	0/6/23/26	0/1/1/1
6	NDG	B	3320	2	-	0/6/23/26	0/1/1/1
4	NAG	B	3371	2	-	0/6/23/26	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(°)	Ideal(°)
4	B	3371	NAG	C2-N2-C7	-2.58	119.72	123.04
6	B	3320	NDG	C2-N2-C7	-2.14	120.29	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2260	NAG	4	0
4	A	2458	NAG	5	0
6	B	3320	NDG	1	0
4	B	3371	NAG	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 [i](#)

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

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

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

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

### 6.3 Carbohydrates [i](#)

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

### 6.4 Ligands [i](#)

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

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

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