



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 06:57 PM GMT

PDB ID : 2VDL
Title : RE-REFINEMENT OF INTEGRIN ALPHAIIIBBETA3 HEADPIECE
Authors : Springer, T.A.; Zhu, J.; Xiao, T.
Deposited on : 2007-10-10
Resolution : 2.75 Å(reported)

This is a full wwPDB validation 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>

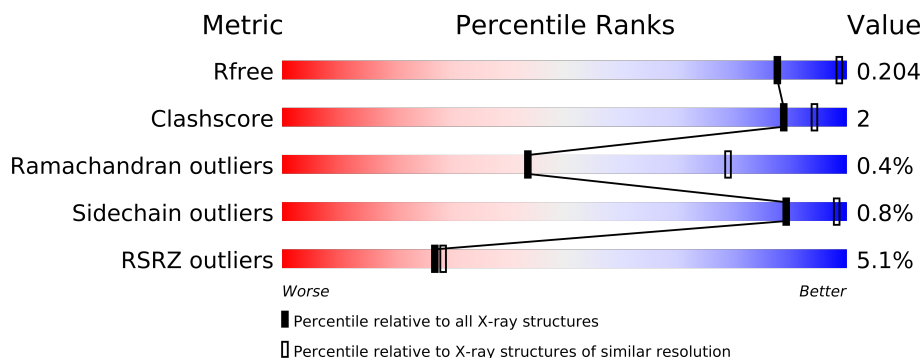
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.75 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	2406 (2.80-2.72)
Clashscore	79885	2995 (2.80-2.72)
Ramachandran outliers	78287	2941 (2.80-2.72)
Sidechain outliers	78261	2944 (2.80-2.72)
RSRZ outliers	66119	2409 (2.80-2.72)

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	A	452	
2	B	461	
3	H	221	
4	L	214	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
5	GOL	A	1453	-	X
5	GOL	B	1463	-	X
7	NAG	B	3099	-	X
8	CAC	B	1462	X	-

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 11800 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 INTEGRIN ALPHA-IIB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	452	Total	C	N	O	S	0	8	0
			3520	2235	612	665	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	282	GLY	ALA	CONFLICT SEE REMARK 9	UNP P08514

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	456	Total	C	N	O	S	0	13	0
			3600	2249	614	704	33			

- Molecule 3 is a protein called MONOCLONAL ANTIBODY 10E5 HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	219	Total	C	N	O	S	0	6	0
			1678	1065	270	337	6			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	214	Total	C	N	O	S	0	7	0
			1663	1035	272	345	11			

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

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

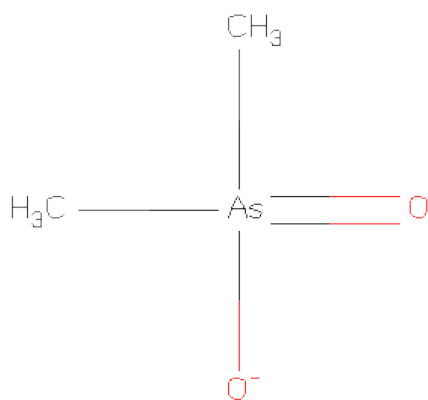
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total	Ca	0	0
			2	2		
6	A	4	Total	Ca	0	0
			4	4		

- Molecule 7 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 8 is CACODYLATE ION (three-letter code: CAC) (formula: $C_2H_6AsO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	B	1	Total	As	C	O	0	0
			5	1	2	2		

- Molecule 9 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 10 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	B	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 11 is a polymer of unknown type called SUGAR (7-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	B	7	Total	C	N	O	0	0
			83	46	2	35		

- Molecule 12 is water.

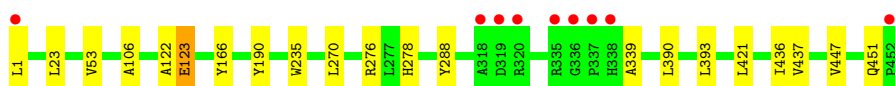
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	411	Total	O	0	0
			411	411		
12	B	250	Total	O	0	0
			250	250		
12	H	221	Total	O	0	0
			221	221		
12	L	247	Total	O	0	0
			247	247		

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.

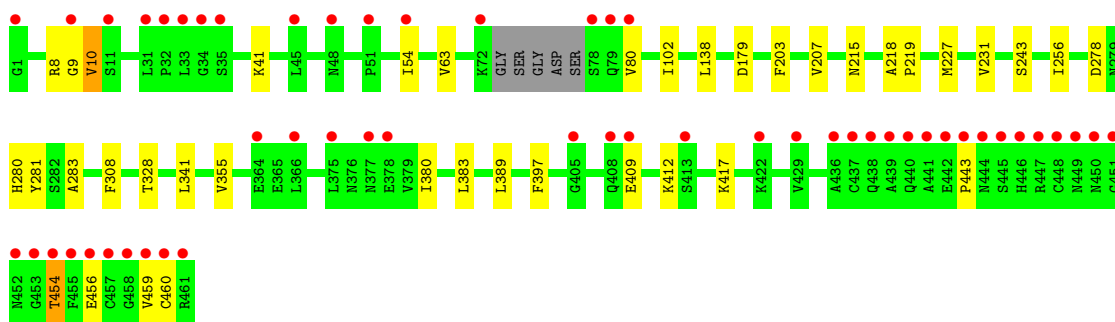
- Molecule 1: INTEGRIN ALPHA-IIB

Chain A: 



- Molecule 2: INTEGRIN BETA-3

Chain B: 



- Molecule 3: MONOCLONAL ANTIBODY 10E5 HEAVY CHAIN

Chain H: 



- Molecule 4: MONOCLONAL ANTIBODY 10E5 LIGHT CHAIN

Chain L: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	148.93Å 148.93Å 176.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.99 – 2.75 38.65 – 2.75	Depositor EDS
% Data completeness (in resolution range)	95.1 (42.99-2.75) 89.6 (38.65-2.75)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.50 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.3.0037	Depositor
R, R_{free}	0.144 , 0.191 0.160 , 0.204	Depositor DCC
R_{free} test set	2696 reflections (5.36%)	DCC
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.005	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.6	EDS
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 53014 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11800	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage'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.*

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: GOL, MG, NAG, CA, CAC, 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	0.34	0/3629	0.52	0/4943
2	B	0.32	0/3688	0.49	0/4995
3	H	0.32	0/1730	0.51	0/2368
4	L	0.34	0/1719	0.53	0/2330
All	All	0.33	0/10766	0.51	0/14636

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
10	B	1	0
11	B	1	0
All	All	2	0

There are no bond length outliers.

There are no bond angle outliers.

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
10	B	3322	MAN	C1
11	B	3373	MAN	C1

There are no planarity outliers.

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	A	3520	0	3361	11	0
2	B	3600	0	3548	20	0
3	H	1678	0	1634	8	0
4	L	1663	0	1581	11	0
5	A	6	0	8	0	0
5	B	6	0	8	0	0
6	A	4	0	0	0	0
6	B	2	0	0	0	0
7	A	28	0	26	0	0
7	B	14	0	13	0	0
8	B	5	0	0	1	0
9	B	1	0	0	0	0
10	B	61	0	52	0	0
11	B	83	0	70	0	0
12	A	411	0	0	2	0
12	B	250	0	0	2	0
12	H	221	0	0	1	0
12	L	247	0	0	2	0
All	All	11800	0	10301	49	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 2.

All (49) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
8:B:1462:CAC:C1	12:B:4099:HOH:O	2.23	0.86
3:H:177[A]:GLN:HB3	4:L:160:LEU:HD11	1.76	0.66
1:A:270:LEU:HD23	1:A:276:ARG:HA	1.80	0.62
2:B:138:LEU:HD23	2:B:341[A]:LEU:HD23	1.84	0.59
1:A:235:TRP:HZ2	1:A:270:LEU:HD11	1.72	0.55
2:B:456:GLU:O	2:B:459:VAL:HG12	2.06	0.55
3:H:177[A]:GLN:NE2	12:H:4196:HOH:O	2.38	0.53
2:B:355:VAL:HG23	2:B:389:LEU:HD22	1.91	0.52
1:A:235:TRP:CZ2	1:A:270:LEU:HD11	2.45	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:390:LEU:HD12	1:A:390:LEU:N	2.25	0.51
1:A:421:LEU:CD2	1:A:437:VAL:HG22	2.40	0.51
2:B:281:TYR:CE1	2:B:283:ALA:HB3	2.46	0.50
1:A:1:LEU:HB2	1:A:393:LEU:HD11	1.94	0.50
2:B:41:LYS:HG3	2:B:54:ILE:HG21	1.93	0.49
4:L:136:LEU:N	4:L:136:LEU:HD12	2.29	0.48
1:A:278[A]:HIS:CD2	1:A:339:ALA:HB1	2.48	0.48
2:B:278:ASP:CG	2:B:280[B]:HIS:HD1	2.17	0.48
1:A:451:GLN:NE2	12:A:4408:HOH:O	2.47	0.48
3:H:172:PHE:CD1	4:L:164:THR:HG23	2.49	0.47
4:L:122:SER:O	4:L:126:THR:HG23	2.16	0.46
2:B:102:ILE:HG22	2:B:397:PHE:HB2	1.97	0.46
2:B:63:VAL:HG23	2:B:63:VAL:O	2.15	0.46
3:H:165:LEU:CD2	3:H:187:VAL:HG21	2.46	0.46
3:H:4:LEU:N	3:H:4:LEU:HD12	2.31	0.46
2:B:308:PHE:CE2	2:B:328:THR:HG21	2.50	0.45
4:L:117:ILE:HG13	4:L:134[B]:CYS:SG	2.57	0.45
3:H:176:LEU:HD13	3:H:181:TYR:CE1	2.51	0.45
2:B:203:PHE:O	2:B:207:VAL:HG23	2.17	0.45
2:B:8:ARG:O	2:B:10:VAL:N	2.51	0.44
3:H:125:PRO:HB3	3:H:151:TYR:HB3	2.00	0.44
4:L:181:LEU:N	4:L:181:LEU:HD23	2.33	0.44
2:B:218:ALA:HB3	2:B:219:PRO:HD3	2.00	0.43
1:A:436:ILE:HG22	1:A:447:VAL:HG22	1.99	0.43
2:B:459:VAL:HG22	2:B:460:CYS:O	2.19	0.43
2:B:278:ASP:OD2	2:B:280[B]:HIS:ND1	2.51	0.43
4:L:105:GLU:HG3	4:L:173:TYR:OH	2.19	0.43
1:A:122:ALA:O	1:A:123:GLU:HB2	2.18	0.42
2:B:227:MET:O	2:B:231:VAL:HG22	2.19	0.42
2:B:380:ILE:HG21	2:B:383:LEU:HD22	2.01	0.42
2:B:417:LYS:NZ	12:B:4224:HOH:O	2.53	0.42
4:L:117:ILE:HA	4:L:134[B]:CYS:SG	2.60	0.42
1:A:53:VAL:HG21	1:A:106:ALA:CB	2.49	0.42
2:B:443:PRO:HA	2:B:454:THR:HG22	2.02	0.42
2:B:409:GLU:HB2	2:B:412[B]:LYS:HZ2	1.84	0.41
4:L:106:ILE:HD12	12:L:4203:HOH:O	2.21	0.41
12:A:4301:HOH:O	2:B:256:ILE:HD12	2.21	0.41
4:L:188[B]:ARG:NH2	12:L:4224:HOH:O	2.54	0.40
3:H:153:PRO:HD2	3:H:207:ALA:CB	2.52	0.40

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	A	457/452 (101%)	445 (97%)	11 (2%)	1 (0%)	56	87
2	B	463/461 (100%)	442 (96%)	17 (4%)	4 (1%)	25	60
3	H	218/221 (99%)	213 (98%)	5 (2%)	0	100	100
4	L	218/214 (102%)	211 (97%)	7 (3%)	0	100	100
All	All	1356/1348 (101%)	1311 (97%)	40 (3%)	5 (0%)	43	79

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	GLU
2	B	9	GLY
2	B	80	VAL
2	B	454	THR
2	B	10	VAL

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	367/360 (102%)	363 (99%)	4 (1%)	84	97
2	B	417/409 (102%)	414 (99%)	3 (1%)	91	98
3	H	192/190 (101%)	191 (100%)	1 (0%)	94	99
4	L	194/188 (103%)	192 (99%)	2 (1%)	85	97
All	All	1170/1147 (102%)	1160 (99%)	10 (1%)	89	97

All (10) 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	179	ASP
2	B	215	ASN
2	B	243	SER
3	H	183	LEU
4	L	194[A]	CYS
4	L	194[B]	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	199	GLN
2	B	316	ASN
2	B	342	GLN
2	B	428	GLN
2	B	438	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

12 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
10	NAG	B	3320	10,2	12,14,15	0.69	1 (8%)	15,19,21	0.95	1 (6%)
10	NAG	B	3321	10	12,14,15	0.76	1 (8%)	15,19,21	1.03	1 (6%)
10	MAN	B	3322	10	10,11,12	0.81	1 (10%)	11,15,17	0.66	0
10	MAN	B	3323	10	10,11,12	0.70	0	11,15,17	0.67	0
10	MAN	B	3324	10	10,11,12	0.70	0	11,15,17	1.09	1 (9%)
11	NAG	B	3371	11,2	12,14,15	0.81	1 (8%)	15,19,21	0.77	0
11	NAG	B	3372	11	12,14,15	0.63	0	15,19,21	0.99	1 (6%)
11	MAN	B	3373	11	10,11,12	0.73	0	11,15,17	1.19	1 (9%)
11	MAN	B	3374	11	10,11,12	0.74	0	11,15,17	1.31	1 (9%)
11	MAN	B	3375	11	10,11,12	0.91	1 (10%)	11,15,17	0.82	0
11	MAN	B	3376	11	10,11,12	0.76	0	11,15,17	1.10	1 (9%)
11	MAN	B	3377	11	10,11,12	0.75	0	11,15,17	0.66	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
10	NAG	B	3320	10,2	-	0/6/23/26	0/1/1/1
10	NAG	B	3321	10	-	0/6/23/26	0/1/1/1
10	MAN	B	3322	10	1/1/4/5	0/2/19/22	0/1/1/1
10	MAN	B	3323	10	-	0/2/19/22	0/1/1/1
10	MAN	B	3324	10	-	0/2/19/22	0/1/1/1
11	NAG	B	3371	11,2	-	0/6/23/26	0/1/1/1
11	NAG	B	3372	11	-	0/6/23/26	0/1/1/1
11	MAN	B	3373	11	1/1/4/5	0/2/19/22	0/1/1/1
11	MAN	B	3374	11	-	0/2/19/22	0/1/1/1
11	MAN	B	3375	11	-	0/2/19/22	0/1/1/1
11	MAN	B	3376	11	-	0/2/19/22	0/1/1/1
11	MAN	B	3377	11	-	0/2/19/22	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	B	3371	NAG	O5-C5	-2.55	1.40	1.45
10	B	3321	NAG	O5-C5	-2.38	1.41	1.45
11	B	3375	MAN	O5-C5	-2.28	1.41	1.45
10	B	3320	NAG	O5-C5	-2.08	1.41	1.45
10	B	3322	MAN	O5-C5	-2.02	1.41	1.45

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	B	3373	MAN	O5-C5-C6	3.33	110.48	106.98
10	B	3324	MAN	O5-C5-C6	2.72	109.84	106.98
11	B	3374	MAN	C3-C4-C5	2.72	115.06	110.20
11	B	3372	NAG	O5-C5-C6	2.37	109.47	106.98
10	B	3321	NAG	O5-C5-C4	-2.18	107.88	110.65
10	B	3320	NAG	C3-C2-N2	-2.16	108.48	111.76
11	B	3376	MAN	O5-C5-C6	2.14	109.22	106.98

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
11	B	3373	MAN	C1
10	B	3322	MAN	C1

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

Of 13 ligands modelled in this entry, 7 are monoatomic - leaving 6 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	$\# Z > 2$
5	GOL	A	1453	-	5,5,5	0.38	0	5,5,5	0.23	0
7	NAG	A	3015	1	12,14,15	0.61	0	15,19,21	1.22	2 (13%)
7	NAG	A	3249	1	12,14,15	0.55	0	15,19,21	1.30	1 (6%)
8	CAC	B	1462	9	4,4,4	3.99	4 (100%)	6,6,6	12.84	3 (50%)
5	GOL	B	1463	6	5,5,5	0.37	0	5,5,5	0.20	0
7	NAG	B	3099	2	12,14,15	0.72	1 (8%)	15,19,21	1.35	2 (13%)

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
5	GOL	A	1453	-	-	0/4/4/4	0/0/0/0
7	NAG	A	3015	1	-	0/6/23/26	0/1/1/1
7	NAG	A	3249	1	-	0/6/23/26	0/1/1/1
8	CAC	B	1462	9	-	0/0/0/0	0/0/0/0
5	GOL	B	1463	6	-	0/4/4/4	0/0/0/0
7	NAG	B	3099	2	-	0/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	1462	CAC	O2-AS	5.45	1.74	1.68
8	B	1462	CAC	AS-C2	3.81	1.97	1.90
8	B	1462	CAC	AS-C1	3.67	1.96	1.90
8	B	1462	CAC	O1-AS	2.44	1.74	1.66
7	B	3099	NAG	O5-C5	-2.11	1.41	1.45

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	1462	CAC	O2-AS-O1	-30.97	105.86	112.54
7	A	3249	NAG	O5-C5-C6	4.34	111.53	106.98
8	B	1462	CAC	O2-AS-C1	4.01	112.43	109.15
8	B	1462	CAC	O2-AS-C2	3.55	112.06	109.15
7	B	3099	NAG	O5-C5-C4	-3.09	106.74	110.65
7	B	3099	NAG	O5-C5-C6	2.94	110.06	106.98
7	A	3015	NAG	C3-C2-N2	-2.83	107.45	111.76
7	A	3015	NAG	O5-C5-C6	2.71	109.83	106.98

There are no chirality outliers.

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	A	452/452 (100%)	0.01	9 (1%) 62 65	22, 41, 59, 79	2 (0%)
2	B	456/461 (98%)	0.49	53 (11%) 5 5	27, 51, 92, 105	2 (0%)
3	H	219/221 (99%)	-0.13	5 (2%) 57 61	25, 40, 62, 102	3 (1%)
4	L	214/214 (100%)	-0.31	1 (0%) 88 91	20, 36, 52, 75	2 (0%)
All	All	1341/1348 (99%)	0.10	68 (5%) 27 28	20, 42, 83, 105	9 (0%)

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	453	GLY	7.7
2	B	451	GLY	7.3
2	B	454	THR	7.3
3	H	221	PRO	6.9
2	B	444	ASN	6.8
2	B	448	CYS	6.8
2	B	443	PRO	6.7
2	B	452	ASN	6.5
2	B	449	ASN	6.3
2	B	446	HIS	6.1
2	B	450	ASN	6.1
3	H	137	THR	5.9
1	A	452	PRO	5.8
4	L	214	CYS	5.7
2	B	460	CYS	5.6
2	B	459	VAL	5.5
1	A	1	LEU	5.5
2	B	461	ARG	5.5
2	B	439	ALA	5.0
2	B	33	LEU	4.8
2	B	78	SER	4.6

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Mol	Chain	Res	Type	RSRZ
2	B	441	ALA	4.6
2	B	455	PHE	4.4
1	A	337[A]	PRO	4.4
2	B	447	ARG	4.0
2	B	442	GLU	4.0
2	B	445	SER	3.9
2	B	79	GLN	3.8
1	A	320	ARG	3.8
2	B	408	GLN	3.8
3	H	220	GLY	3.7
2	B	458	GLY	3.7
2	B	440	GLN	3.6
1	A	319	ASP	3.6
2	B	80	VAL	3.5
2	B	1	GLY	3.3
2	B	457	CYS	3.3
1	A	338	HIS	3.1
2	B	456	GLU	3.1
2	B	409	GLU	3.0
2	B	54	ILE	2.9
2	B	48	ASN	2.9
3	H	138	THR	2.9
2	B	34	GLY	2.8
2	B	377	ASN	2.8
2	B	378	GLU	2.8
2	B	437	CYS	2.6
2	B	405	GLY	2.5
2	B	11	SER	2.4
2	B	32	PRO	2.4
1	A	318	ALA	2.4
2	B	364	GLU	2.4
2	B	438	GLN	2.4
2	B	422	LYS	2.4
2	B	35	SER	2.4
2	B	366	LEU	2.4
2	B	436	ALA	2.3
1	A	335[A]	ARG	2.3
1	A	336	GLY	2.3
3	H	211	LYS	2.3
2	B	9	GLY	2.2
2	B	72	LYS	2.2
2	B	51	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	413	SER	2.1
2	B	31	LEU	2.1
2	B	45	LEU	2.1
2	B	375	LEU	2.0
2	B	429	VAL	2.0

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

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

6.3 Carbohydrates ⓘ

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
11	MAN	B	3373	11/12	0.17	3.84	75,86,88,91	0
11	MAN	B	3374	11/12	0.27	2.00	90,94,95,95	0
10	NAG	B	3321	14/15	0.20	1.06	66,69,76,84	0
11	NAG	B	3371	14/15	0.17	0.06	62,67,71,72	0
11	MAN	B	3375	11/12	0.14	0.02	63,70,74,78	0
11	MAN	B	3376	11/12	0.15	-0.08	53,57,61,62	0
10	NAG	B	3320	14/15	0.10	-1.80	41,48,54,60	0
10	MAN	B	3324	11/12	0.36	-	104,106,106,106	0
10	MAN	B	3322	11/12	0.31	-	90,96,101,105	0
11	NAG	B	3372	14/15	0.20	-	56,74,80,84	0
11	MAN	B	3377	11/12	0.29	-	82,85,87,89	0
10	MAN	B	3323	11/12	0.29	-	91,94,96,96	0

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
7	NAG	B	3099	14/15	0.28	4.10	94,100,104,104	0
5	GOL	B	1463	6/6	0.18	3.63	57,58,60,61	0
5	GOL	A	1453	6/6	0.29	2.21	58,62,65,66	0
8	CAC	B	1462	5/5	0.13	-0.97	51,52,53,55	0
6	CA	A	2004	1/1	0.08	-1.37	57,57,57,57	0
6	CA	A	2006	1/1	0.10	-1.56	57,57,57,57	0
6	CA	A	2005	1/1	0.06	-2.23	41,41,41,41	0
6	CA	A	2007	1/1	0.06	-2.42	53,53,53,53	0
6	CA	B	2002	1/1	0.06	-2.88	32,32,32,32	0
9	MG	B	2001	1/1	0.07	-3.01	23,23,23,23	0
6	CA	B	2003	1/1	0.07	-4.42	25,25,25,25	0
7	NAG	A	3015	14/15	0.39	-	73,74,75,76	14
7	NAG	A	3249	14/15	0.50	-	85,87,88,89	14

6.5 Other polymers ⓘ

There are no such residues in this entry.