



Full wwPDB X-ray Structure Validation Report

Nov 6, 2014 – 01:10 PM EST

PDB ID : 4CDH
Title : Crystallographic structure of the Human Iggl alpha 2-6 sialilated Fc-Fragment
Authors : Silva-Martin, N.; Bartual, S.G.; Hermoso, J.A.
Deposited on : 2013-10-31
Resolution : 2.30 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>

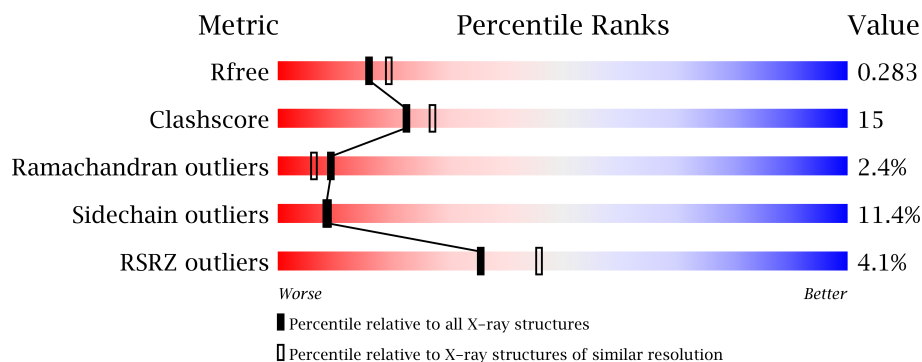
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable24103
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.1.3
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable24103

1 Overall quality at a glance

The reported resolution of this entry is 2.30 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	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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	255	
1	B	255	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3619 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 IG GAMMA-1 CHAIN C REGION.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	S	0	0	0
			1654	1054	278	316	6			
1	B	208	Total	C	N	O	S	0	0	0
			1662	1058	280	318	6			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-26	MET	-	EXPRESSION TAG	UNP P01857
A	-25	LYS	-	EXPRESSION TAG	UNP P01857
A	-24	LYS	-	EXPRESSION TAG	UNP P01857
A	-23	THR	-	EXPRESSION TAG	UNP P01857
A	-22	ALA	-	EXPRESSION TAG	UNP P01857
A	-21	ILE	-	EXPRESSION TAG	UNP P01857
A	-20	ALA	-	EXPRESSION TAG	UNP P01857
A	-19	ILE	-	EXPRESSION TAG	UNP P01857
A	-18	ALA	-	EXPRESSION TAG	UNP P01857
A	-17	VAL	-	EXPRESSION TAG	UNP P01857
A	-16	ALA	-	EXPRESSION TAG	UNP P01857
A	-15	LEU	-	EXPRESSION TAG	UNP P01857
A	-14	ALA	-	EXPRESSION TAG	UNP P01857
A	-13	GLY	-	EXPRESSION TAG	UNP P01857
A	-12	PHE	-	EXPRESSION TAG	UNP P01857
A	-11	ALA	-	EXPRESSION TAG	UNP P01857
A	-10	THR	-	EXPRESSION TAG	UNP P01857
A	-9	VAL	-	EXPRESSION TAG	UNP P01857
A	-8	ALA	-	EXPRESSION TAG	UNP P01857
A	-7	GLN	-	EXPRESSION TAG	UNP P01857
A	-6	ALA	-	EXPRESSION TAG	UNP P01857
A	-5	ASP	-	EXPRESSION TAG	UNP P01857
A	-4	VAL	-	EXPRESSION TAG	UNP P01857
A	-3	GLU	-	EXPRESSION TAG	UNP P01857
A	-2	SER	-	EXPRESSION TAG	UNP P01857

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-26	MET	-	EXPRESSION TAG	UNP P01857
B	-25	LYS	-	EXPRESSION TAG	UNP P01857
B	-24	LYS	-	EXPRESSION TAG	UNP P01857
B	-23	THR	-	EXPRESSION TAG	UNP P01857
B	-22	ALA	-	EXPRESSION TAG	UNP P01857
B	-21	ILE	-	EXPRESSION TAG	UNP P01857
B	-20	ALA	-	EXPRESSION TAG	UNP P01857
B	-19	ILE	-	EXPRESSION TAG	UNP P01857
B	-18	ALA	-	EXPRESSION TAG	UNP P01857
B	-17	VAL	-	EXPRESSION TAG	UNP P01857
B	-16	ALA	-	EXPRESSION TAG	UNP P01857
B	-15	LEU	-	EXPRESSION TAG	UNP P01857
B	-14	ALA	-	EXPRESSION TAG	UNP P01857
B	-13	GLY	-	EXPRESSION TAG	UNP P01857
B	-12	PHE	-	EXPRESSION TAG	UNP P01857
B	-11	ALA	-	EXPRESSION TAG	UNP P01857
B	-10	THR	-	EXPRESSION TAG	UNP P01857
B	-9	VAL	-	EXPRESSION TAG	UNP P01857
B	-8	ALA	-	EXPRESSION TAG	UNP P01857
B	-7	GLN	-	EXPRESSION TAG	UNP P01857
B	-6	ALA	-	EXPRESSION TAG	UNP P01857
B	-5	ASP	-	EXPRESSION TAG	UNP P01857
B	-4	VAL	-	EXPRESSION TAG	UNP P01857
B	-3	GLU	-	EXPRESSION TAG	UNP P01857
B	-2	SER	-	EXPRESSION TAG	UNP P01857

- Molecule 2 is a polymer of unknown type called SUGAR (10-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	10	Total	C	N	O	0	0
			121	68	4	49		

- Molecule 3 is a polymer of unknown type called SUGAR (8-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	8	Total	C	N	O	0	0
			100	56	4	40		

- Molecule 4 is water.

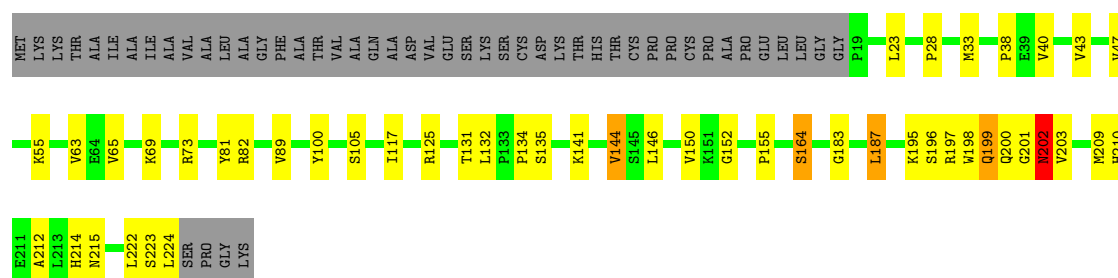
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	49	Total 49	O 49	0	0
4	B	33	Total 33	O 33	0	0

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.

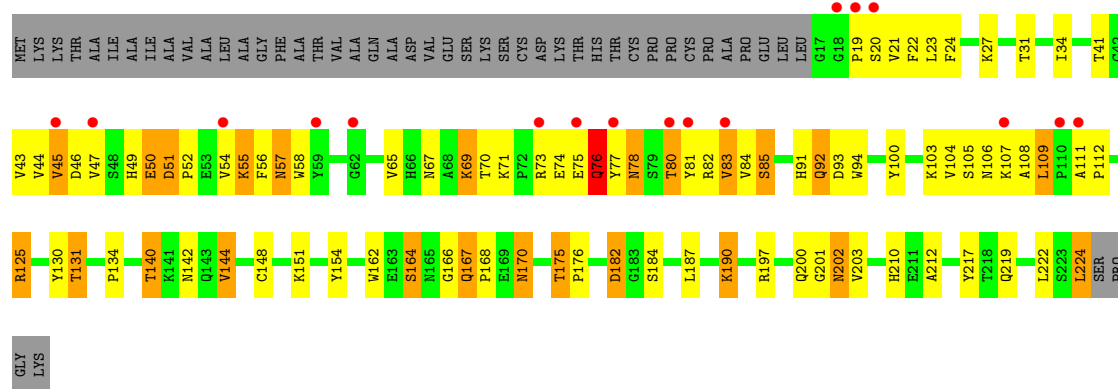
• Molecule 1: IG GAMMA-1 CHAIN C REGION

Chain A: 



• Molecule 1: IG GAMMA-1 CHAIN C REGION

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.59Å 80.15Å 140.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.10 – 2.30 42.17 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.10-2.30) 100.0 (42.17-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.234 , 0.282 0.237 , 0.283	Depositor DCC
R_{free} test set	1300 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	38.3	Xtriage
Anisotropy	0.112	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 25579 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3619	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.97% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: GAL, FUL, BMA, 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	0.77	0/1700	0.86	0/2316
1	B	0.81	0/1708	0.93	1/2327 (0.0%)
All	All	0.79	0/3408	0.90	1/4643 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	182	ASP	CB-CA-C	-5.78	98.84	110.40

There are no chirality outliers.

There are no planarity outliers.

5.2 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 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	1654	0	1623	35	0
1	B	1662	0	1628	68	0
2	A	121	0	103	1	0
3	B	100	0	85	4	0
4	A	49	0	0	2	0
4	B	33	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3619	0	3439	103	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.

All (103) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:41:THR:HG21	3:B:1005:NAG:O6	1.67	0.94
1:B:131:THR:HG23	1:B:222:LEU:HD13	1.60	0.82
1:B:201:GLY:HA2	1:B:224:LEU:HD11	1.61	0.81
1:B:76:GLN:HE22	1:B:82:ARG:HG3	1.46	0.81
1:B:91:HIS:HA	4:B:2010:HOH:O	1.80	0.80
1:A:164:SER:OG	1:A:203:VAL:O	2.01	0.78
1:B:75:GLU:O	1:B:76:GLN:HB2	1.85	0.76
1:B:164:SER:OG	1:B:203:VAL:O	2.04	0.75
1:B:92:GLN:N	4:B:2010:HOH:O	2.15	0.73
1:A:198:TRP:O	1:A:224:LEU:HD11	1.89	0.72
1:A:33:MET:CE	1:A:209:MET:SD	2.77	0.72
1:B:41:THR:HG21	3:B:1005:NAG:HO6	1.55	0.71
1:A:125:ARG:HG3	1:A:152:GLY:O	1.89	0.71
1:B:210:HIS:HD2	1:B:212:ALA:H	1.39	0.70
1:A:198:TRP:O	1:A:224:LEU:CD1	2.39	0.70
1:A:195:LYS:O	1:A:199:GLN:HG2	1.93	0.68
1:B:20:SER:O	1:B:44:VAL:O	2.10	0.68
1:B:81:TYR:HB3	4:B:2008:HOH:O	1.92	0.68
1:A:33:MET:HE2	1:A:209:MET:SD	2.36	0.65
1:A:33:MET:HE3	1:A:209:MET:SD	2.38	0.63
1:A:125:ARG:CG	1:A:152:GLY:O	2.47	0.63
1:A:134:PRO:HG3	1:A:144:VAL:HG13	1.82	0.61
1:A:150:VAL:HB	1:A:187:LEU:CD2	2.32	0.60
1:B:41:THR:CG2	3:B:1005:NAG:O6	2.46	0.60
1:B:224:LEU:H	1:B:224:LEU:HD12	1.67	0.60
1:A:135:SER:HB2	1:B:130:TYR:HB3	1.85	0.59
1:B:140:THR:OG1	1:B:140:THR:O	2.20	0.59
1:A:210:HIS:CD2	1:A:212:ALA:H	2.21	0.58
1:B:182:ASP:HB3	1:B:184:SER:H	1.68	0.58
1:B:21:VAL:O	1:B:22:PHE:CD1	2.56	0.58
1:A:201:GLY:CA	1:A:202:ASN:HB2	2.33	0.58
1:B:190:LYS:HD2	4:B:2024:HOH:O	2.03	0.58
1:A:89:VAL:HG22	1:A:100:TYR:CZ	2.40	0.57
1:A:134:PRO:HG3	1:A:144:VAL:CG1	2.37	0.55
1:B:19:PRO:HB3	1:B:46:ASP:O	2.06	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:89:VAL:HG22	1:A:100:TYR:CE2	2.43	0.54
1:A:155:PRO:O	1:A:210:HIS:HE1	1.90	0.54
1:A:82:ARG:NH1	2:A:1002:NAG:O7	2.40	0.54
4:A:2033:HOH:O	1:B:190:LYS:HD3	2.07	0.54
1:B:125:ARG:HG3	1:B:154:TYR:HB3	1.90	0.54
1:A:183:GLY:HA2	4:A:2043:HOH:O	2.07	0.53
1:B:91:HIS:CA	4:B:2010:HOH:O	2.48	0.53
1:B:47:VAL:O	1:B:80:THR:HG22	2.09	0.52
1:B:71:LYS:HB3	1:B:84:VAL:HG22	1.90	0.52
1:B:45:VAL:HG12	1:B:46:ASP:H	1.75	0.52
1:B:175:THR:HG23	1:B:176:PRO:O	2.11	0.51
1:A:131:THR:HG21	1:A:222:LEU:HB2	1.92	0.51
1:B:109:LEU:HD22	1:B:111:ALA:O	2.11	0.51
1:A:201:GLY:HA3	1:A:202:ASN:HB2	1.92	0.51
1:A:23:LEU:HD11	1:A:40:VAL:CG1	2.41	0.51
1:A:131:THR:HG22	1:A:222:LEU:HD22	1.94	0.50
1:B:47:VAL:H	1:B:80:THR:HB	1.76	0.49
1:A:55:LYS:HB3	1:A:105:SER:HB2	1.92	0.49
1:B:77:TYR:O	1:B:78:ASN:HB3	2.12	0.49
1:A:43:VAL:HG11	1:A:82:ARG:NH1	2.28	0.49
1:B:91:HIS:O	1:B:92:GLN:HB2	2.13	0.49
1:B:210:HIS:CD2	1:B:212:ALA:H	2.25	0.48
1:A:131:THR:O	1:A:132:LEU:HD12	2.14	0.48
3:B:1001:NAG:H82	3:B:1001:NAG:O3	2.15	0.47
1:B:47:VAL:N	1:B:80:THR:HB	2.29	0.47
1:B:111:ALA:HB1	1:B:112:PRO:HD2	1.97	0.47
1:B:70:THR:HA	1:B:85:SER:HA	1.97	0.47
1:A:210:HIS:HD2	1:A:212:ALA:H	1.62	0.46
1:B:56:PHE:HZ	1:B:83:VAL:O	1.99	0.46
1:B:106:ASN:OD1	1:B:108:ALA:N	2.47	0.46
1:B:58:TRP:HE1	1:B:85:SER:HG	1.64	0.45
1:B:71:LYS:HB3	1:B:84:VAL:CG2	2.46	0.45
1:B:24:PHE:CD1	1:B:24:PHE:N	2.85	0.45
1:A:134:PRO:HD3	1:A:146:LEU:HD23	2.00	0.44
1:B:134:PRO:HB3	1:B:144:VAL:HG22	1.98	0.44
1:B:202:ASN:N	1:B:202:ASN:OD1	2.51	0.44
1:B:91:HIS:CD2	1:B:91:HIS:H	2.35	0.44
1:B:170:ASN:OD1	1:B:170:ASN:C	2.55	0.44
1:B:51:ASP:OD1	1:B:108:ALA:HB2	2.18	0.44
1:B:69:LYS:HA	1:B:69:LYS:CE	2.48	0.43
1:A:38:PRO:HG2	1:A:89:VAL:O	2.18	0.43
1:B:54:VAL:HG12	1:B:55:LYS:N	2.33	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:201:GLY:CA	1:B:224:LEU:HD11	2.40	0.43
1:B:74:GLU:O	1:B:81:TYR:HB2	2.19	0.43
1:A:201:GLY:CA	1:A:202:ASN:CB	2.97	0.43
1:B:57:ASN:HB2	1:B:103:LYS:HB3	2.00	0.43
1:B:166:GLY:C	1:B:167:GLN:HE21	2.22	0.43
1:B:175:THR:HG22	1:B:187:LEU:HB2	2.01	0.43
1:B:45:VAL:HG12	1:B:46:ASP:N	2.33	0.43
1:A:214:HIS:O	1:A:215:ASN:HB2	2.18	0.43
1:B:187:LEU:C	1:B:187:LEU:HD12	2.39	0.42
1:B:50:GLU:C	1:B:52:PRO:HD3	2.39	0.42
1:B:77:TYR:O	1:B:78:ASN:CB	2.67	0.42
1:B:104:VAL:HG12	1:B:105:SER:N	2.35	0.42
1:B:167:GLN:HA	1:B:168:PRO:HD3	1.89	0.42
1:B:69:LYS:HE2	1:B:69:LYS:HA	2.02	0.42
1:A:131:THR:C	1:A:132:LEU:HD12	2.39	0.41
1:B:55:LYS:HB3	1:B:105:SER:HB2	2.02	0.41
1:B:182:ASP:CB	1:B:184:SER:H	2.33	0.41
1:B:148:CYS:HB2	1:B:162:TRP:CZ2	2.55	0.41
1:B:175:THR:CG2	1:B:176:PRO:O	2.68	0.41
1:B:217:TYR:CE1	1:B:219:GLN:HG3	2.55	0.41
1:B:71:LYS:HD3	1:B:84:VAL:HG21	2.03	0.41
1:B:93:ASP:HB3	1:B:100:TYR:OH	2.21	0.41
1:B:31:THR:HG21	1:B:94:TRP:CD1	2.56	0.41
1:A:134:PRO:HD3	1:A:146:LEU:CD2	2.51	0.40
1:A:47:VAL:HB	1:A:81:TYR:HB2	2.02	0.40
1:B:200:GLN:CG	1:B:202:ASN:OD1	2.70	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	204/255 (80%)	197 (97%)	6 (3%)	1 (0%)	38	45

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	206/255 (81%)	176 (85%)	21 (10%)	9 (4%)	4	1
All	All	410/510 (80%)	373 (91%)	27 (7%)	10 (2%)	9	6

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	45	VAL
1	B	78	ASN
1	B	76	GLN
1	B	80	THR
1	B	85	SER
1	B	92	GLN
1	A	202	ASN
1	B	67	ASN
1	B	50	GLU
1	B	83	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	193/229 (84%)	177 (92%)	16 (8%)	16	19
1	B	193/229 (84%)	165 (86%)	28 (14%)	5	4
All	All	386/458 (84%)	342 (89%)	44 (11%)	8	9

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	PRO
1	A	63	VAL
1	A	65	VAL
1	A	69	LYS
1	A	73	ARG
1	A	117	ILE
1	A	141	LYS
1	A	144	VAL

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Mol	Chain	Res	Type
1	A	164	SER
1	A	187	LEU
1	A	196	SER
1	A	197	ARG
1	A	199	GLN
1	A	200	GLN
1	A	202	ASN
1	A	223	SER
1	B	23	LEU
1	B	27	LYS
1	B	34	ILE
1	B	43	VAL
1	B	49	HIS
1	B	51	ASP
1	B	55	LYS
1	B	57	ASN
1	B	65	VAL
1	B	69	LYS
1	B	73	ARG
1	B	76	GLN
1	B	107	LYS
1	B	109	LEU
1	B	125	ARG
1	B	131	THR
1	B	140	THR
1	B	142	ASN
1	B	144	VAL
1	B	151	LYS
1	B	164	SER
1	B	167	GLN
1	B	170	ASN
1	B	175	THR
1	B	190	LYS
1	B	197	ARG
1	B	202	ASN
1	B	224	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS
1	A	76	GLN

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Mol	Chain	Res	Type
1	A	92	GLN
1	A	199	GLN
1	A	202	ASN
1	A	210	HIS
1	A	219	GLN
1	B	76	GLN
1	B	91	HIS
1	B	142	ASN
1	B	167	GLN
1	B	210	HIS

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 ⓘ

18 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	$\# Z > 2$
2	NAG	A	1001	1,2	12,14,15	0.61	0	15,19,21	2.02	5 (33%)
2	NAG	A	1002	2	12,14,15	1.25	2 (16%)	15,19,21	1.37	2 (13%)
2	BMA	A	1003	2	10,11,12	1.10	1 (10%)	11,15,17	1.66	3 (27%)
2	MAN	A	1004	2	10,11,12	1.01	1 (10%)	11,15,17	2.96	6 (54%)
2	NAG	A	1005	2	12,14,15	1.20	1 (8%)	15,19,21	2.04	4 (26%)
2	GAL	A	1006	2	10,11,12	1.00	0	11,15,17	2.34	4 (36%)
2	MAN	A	1008	2	10,11,12	0.86	1 (10%)	11,15,17	2.15	4 (36%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	A	1009	2	12,14,15	0.83	0	15,19,21	2.36	6 (40%)
2	GAL	A	1010	2	10,11,12	0.90	0	11,15,17	1.76	3 (27%)
2	FUL	A	1011	2	9,10,11	1.11	1 (11%)	10,14,16	0.89	0
3	NAG	B	1001	1,3	12,14,15	0.61	0	15,19,21	2.14	6 (40%)
3	NAG	B	1002	3	12,14,15	0.76	0	15,19,21	1.16	2 (13%)
3	BMA	B	1003	3	10,11,12	0.91	0	11,15,17	1.61	3 (27%)
3	MAN	B	1004	3	10,11,12	0.88	0	11,15,17	1.89	3 (27%)
3	NAG	B	1005	3	12,14,15	0.92	1 (8%)	15,19,21	2.50	8 (53%)
3	GAL	B	1006	3	10,11,12	0.90	0	11,15,17	3.59	4 (36%)
3	MAN	B	1008	3	10,11,12	0.73	0	11,15,17	2.59	4 (36%)
3	NAG	B	1009	3	12,14,15	0.77	0	15,19,21	2.42	6 (40%)

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
2	NAG	A	1001	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1002	2	-	0/6/23/26	0/1/1/1
2	BMA	A	1003	2	-	0/2/19/22	0/1/1/1
2	MAN	A	1004	2	-	0/2/19/22	0/1/1/1
2	NAG	A	1005	2	-	0/6/23/26	0/1/1/1
2	GAL	A	1006	2	-	0/2/19/22	0/1/1/1
2	MAN	A	1008	2	-	0/2/19/22	0/1/1/1
2	NAG	A	1009	2	-	0/6/23/26	0/1/1/1
2	GAL	A	1010	2	-	0/2/19/22	0/1/1/1
2	FUL	A	1011	2	-	0/0/17/20	0/1/1/1
3	NAG	B	1001	1,3	-	0/6/23/26	0/1/1/1
3	NAG	B	1002	3	-	0/6/23/26	0/1/1/1
3	BMA	B	1003	3	-	0/2/19/22	0/1/1/1
3	MAN	B	1004	3	-	0/2/19/22	0/1/1/1
3	NAG	B	1005	3	-	0/6/23/26	0/1/1/1
3	GAL	B	1006	3	-	0/2/19/22	0/1/1/1
3	MAN	B	1008	3	-	0/2/19/22	0/1/1/1
3	NAG	B	1009	3	-	0/6/23/26	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1005	NAG	O5-C5	-3.54	1.39	1.45
2	A	1002	NAG	C2-N2	-3.01	1.43	1.46
2	A	1002	NAG	O5-C5	-2.70	1.41	1.45
2	A	1011	FUL	O5-C5	-2.63	1.41	1.45
2	A	1003	BMA	O5-C5	-2.56	1.41	1.45
3	B	1005	NAG	O5-C5	-2.32	1.41	1.45
2	A	1008	MAN	O5-C5	-2.20	1.42	1.45
2	A	1004	MAN	O5-C5	-2.17	1.42	1.45

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1006	GAL	O5-C5-C6	9.08	116.50	106.98
2	A	1004	MAN	O5-C5-C4	-7.44	101.21	110.65
3	B	1008	MAN	O5-C5-C6	7.06	114.39	106.98
3	B	1009	NAG	O5-C5-C6	6.43	113.72	106.98
2	A	1006	GAL	O5-C5-C6	5.43	112.68	106.98
3	B	1006	GAL	O5-C5-C4	-5.35	103.86	110.65
3	B	1005	NAG	O5-C5-C4	4.53	116.40	110.65
2	A	1009	NAG	C3-C2-N2	4.48	118.34	111.62
2	A	1005	NAG	C8-C7-N2	4.47	124.57	116.12
2	A	1009	NAG	C2-N2-C7	4.43	128.75	123.39
3	B	1004	MAN	C4-C3-C2	-4.32	105.79	110.61
2	A	1010	GAL	C3-C4-C5	4.03	117.41	110.17
2	A	1008	MAN	O2-C2-C3	-4.03	101.56	110.10
3	B	1005	NAG	C3-C2-N2	3.93	117.53	111.62
2	A	1001	NAG	O5-C5-C4	3.91	115.62	110.65
3	B	1001	NAG	C8-C7-N2	3.82	123.34	116.12
3	B	1001	NAG	O5-C5-C6	3.62	110.78	106.98
3	B	1006	GAL	C4-C3-C2	-3.61	106.59	110.61
2	A	1001	NAG	C3-C2-N2	3.59	117.02	111.62
2	A	1008	MAN	C4-C3-C2	-3.50	106.70	110.61
2	A	1005	NAG	O5-C5-C6	-3.39	103.42	106.98
3	B	1005	NAG	C8-C7-N2	3.30	122.36	116.12
2	A	1006	GAL	C4-C3-C2	-3.27	106.96	110.61
2	A	1009	NAG	O3-C3-C4	3.27	117.65	110.36
3	B	1001	NAG	C3-C4-C5	3.24	115.98	110.17
3	B	1005	NAG	O7-C7-C8	-3.23	115.84	122.04
3	B	1009	NAG	C3-C4-C5	3.23	115.97	110.17
3	B	1009	NAG	C8-C7-N2	3.19	122.15	116.12
2	A	1004	MAN	C6-C5-C4	3.16	120.76	113.04
3	B	1009	NAG	C4-C3-C2	3.14	117.34	110.74
3	B	1006	GAL	O2-C2-C3	-3.10	103.53	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1009	NAG	O5-C5-C6	3.09	110.22	106.98
2	A	1003	BMA	O5-C5-C6	3.03	110.16	106.98
2	A	1002	NAG	C3-C2-N2	-3.02	107.10	111.62
3	B	1008	MAN	C3-C4-C5	-2.97	104.83	110.17
3	B	1005	NAG	O3-C3-C4	-2.94	103.78	110.36
2	A	1001	NAG	C2-N2-C7	2.94	126.95	123.39
2	A	1004	MAN	C4-C3-C2	-2.90	107.37	110.61
2	A	1006	GAL	C6-C5-C4	-2.90	105.95	113.04
2	A	1010	GAL	O5-C5-C4	2.90	114.33	110.65
2	A	1008	MAN	O4-C4-C5	2.89	116.90	109.25
2	A	1006	GAL	O2-C2-C3	-2.81	104.15	110.10
3	B	1005	NAG	O5-C5-C6	-2.79	104.05	106.98
3	B	1001	NAG	C2-N2-C7	2.77	126.75	123.39
2	A	1001	NAG	O5-C5-C6	2.76	109.88	106.98
3	B	1001	NAG	O7-C7-C8	-2.69	116.89	122.04
2	A	1004	MAN	O5-C5-C6	2.62	109.73	106.98
2	A	1005	NAG	O7-C7-C8	-2.59	117.07	122.04
3	B	1005	NAG	C3-C4-C5	2.57	114.78	110.17
2	A	1002	NAG	C3-C4-C5	-2.57	105.56	110.17
2	A	1004	MAN	O2-C2-C3	-2.52	104.75	110.10
3	B	1003	BMA	O3-C3-C4	-2.51	104.75	110.36
3	B	1009	NAG	O5-C5-C4	2.45	113.76	110.65
3	B	1002	NAG	O5-C5-C6	2.43	109.53	106.98
3	B	1002	NAG	O4-C4-C5	2.41	115.63	109.25
2	A	1003	BMA	O4-C4-C3	-2.38	105.04	110.36
3	B	1005	NAG	O3-C3-C2	2.35	113.97	109.16
3	B	1008	MAN	O3-C3-C2	2.31	113.94	109.74
2	A	1008	MAN	O4-C4-C3	-2.29	105.23	110.36
2	A	1009	NAG	O4-C4-C5	2.29	115.32	109.25
3	B	1008	MAN	C4-C3-C2	-2.29	108.06	110.61
3	B	1004	MAN	O4-C4-C3	-2.27	105.29	110.36
2	A	1001	NAG	C6-C5-C4	-2.26	107.51	113.04
2	A	1005	NAG	C3-C2-N2	2.26	115.02	111.62
3	B	1003	BMA	O5-C5-C6	2.24	109.33	106.98
3	B	1004	MAN	O5-C5-C4	2.09	113.31	110.65
2	A	1003	BMA	O3-C3-C4	-2.08	105.71	110.36
2	A	1009	NAG	O4-C4-C3	2.07	114.98	110.36
3	B	1001	NAG	C3-C2-N2	-2.07	108.52	111.62
3	B	1003	BMA	C3-C4-C5	2.07	113.88	110.17
2	A	1004	MAN	O4-C4-C5	2.04	114.66	109.25
3	B	1009	NAG	O7-C7-C8	-2.04	118.14	122.04
2	A	1010	GAL	O5-C5-C6	2.02	109.10	106.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	206/255 (80%)	-0.19	0 100 100	21, 35, 58, 69	0
1	B	208/255 (81%)	0.42	17 (8%) 12 18	18, 45, 82, 98	0
All	All	414/510 (81%)	0.11	17 (4%) 35 46	18, 39, 77, 98	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	83	VAL	5.9
1	B	81	TYR	5.1
1	B	20	SER	4.2
1	B	54	VAL	3.7
1	B	77	TYR	3.7
1	B	18	GLY	3.6
1	B	111	ALA	2.7
1	B	62	GLY	2.6
1	B	107	LYS	2.6
1	B	45	VAL	2.5
1	B	19	PRO	2.5
1	B	75	GLU	2.3
1	B	47	VAL	2.2
1	B	80	THR	2.2
1	B	59	TYR	2.2
1	B	110	PRO	2.0
1	B	73	ARG	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
2	FUL	A	1011	10/11	0.21	5.50	63,69,72,76	0
2	GAL	A	1010	11/12	0.18	1.32	69,76,84,85	0
3	NAG	B	1001	14/15	0.24	0.30	60,74,86,92	0
2	GAL	A	1006	11/12	0.12	-0.05	30,36,42,51	0
2	NAG	A	1005	14/15	0.11	-0.18	40,46,52,58	0
2	BMA	A	1003	11/12	0.12	-0.33	37,44,53,54	0
3	GAL	B	1006	11/12	0.12	-0.50	40,43,52,53	0
2	NAG	A	1002	14/15	0.12	-0.61	34,38,43,46	0
3	NAG	B	1002	14/15	0.20	-0.77	68,76,87,90	0
3	NAG	B	1005	14/15	0.13	-1.20	50,56,66,70	0
3	MAN	B	1004	11/12	0.10	-1.95	59,62,64,65	0
3	BMA	B	1003	11/12	0.12	-2.03	61,67,71,74	0
2	NAG	A	1001	14/15	0.10	-2.80	42,49,63,66	0
2	MAN	A	1004	11/12	0.08	-2.95	41,47,50,55	0
3	NAG	B	1009	14/15	0.19	-	69,78,87,87	0
2	NAG	A	1009	14/15	0.21	-	56,77,84,86	0
3	MAN	B	1008	11/12	0.14	-	63,67,69,75	0
2	MAN	A	1008	11/12	0.14	-	60,63,68,68	0

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.