



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:15 PM GMT

PDB ID : 4BYH
Title : Crystal structure of sialylated IgG Fc
Authors : Crispin, M.; Yu, X.; Bowden, T.A.
Deposited on : 2013-07-19
Resolution : 2.30 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

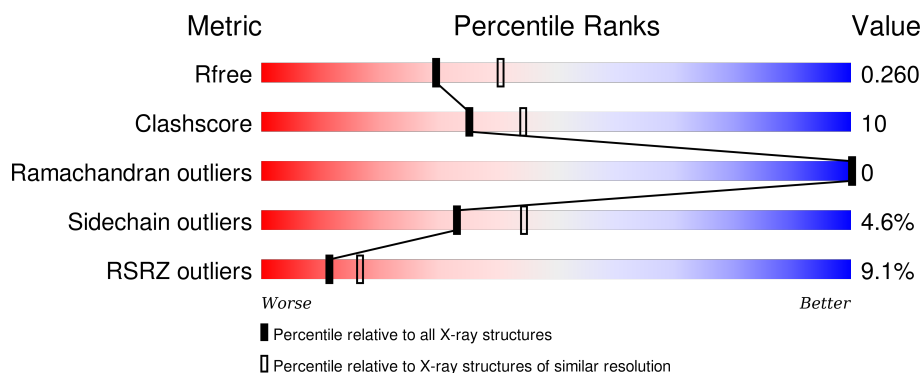
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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. 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.

Mol	Chain	Length	Quality of chain
1	A	233	<div> <div>15%</div> <div>75%</div> <div>12%</div> <div>•</div> <div>11%</div> </div>
1	B	233	<div> <div>%</div> <div>79%</div> <div>10%</div> <div>10%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3780 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 IG GAMMA-1 CHAIN C REGION.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1661	1057	279	318	7			
1	B	209	Total	C	N	O	S	0	5	0
			1700	1083	286	324	7			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	222	GLY	-	EXPRESSION TAG	UNP P01857
A	356	GLU	ASP	CONFLICT	UNP P01857
A	358	MET	LEU	CONFLICT	UNP P01857
A	447	THR	-	EXPRESSION TAG	UNP P01857
A	448	LYS	-	EXPRESSION TAG	UNP P01857
A	449	HIS	-	EXPRESSION TAG	UNP P01857
A	450	HIS	-	EXPRESSION TAG	UNP P01857
A	451	HIS	-	EXPRESSION TAG	UNP P01857
A	452	HIS	-	EXPRESSION TAG	UNP P01857
A	453	HIS	-	EXPRESSION TAG	UNP P01857
A	454	HIS	-	EXPRESSION TAG	UNP P01857
B	222	GLY	-	EXPRESSION TAG	UNP P01857
B	356	GLU	ASP	CONFLICT	UNP P01857
B	358	MET	LEU	CONFLICT	UNP P01857
B	447	THR	-	EXPRESSION TAG	UNP P01857
B	448	LYS	-	EXPRESSION TAG	UNP P01857
B	449	HIS	-	EXPRESSION TAG	UNP P01857
B	450	HIS	-	EXPRESSION TAG	UNP P01857
B	451	HIS	-	EXPRESSION TAG	UNP P01857
B	452	HIS	-	EXPRESSION TAG	UNP P01857
B	453	HIS	-	EXPRESSION TAG	UNP P01857
B	454	HIS	-	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
			130	73	5	52		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		
3	A	1	Total	Cl	0	0
			1	1		

- Molecule 4 is a polymer of unknown type called SUGAR (10-MER).

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

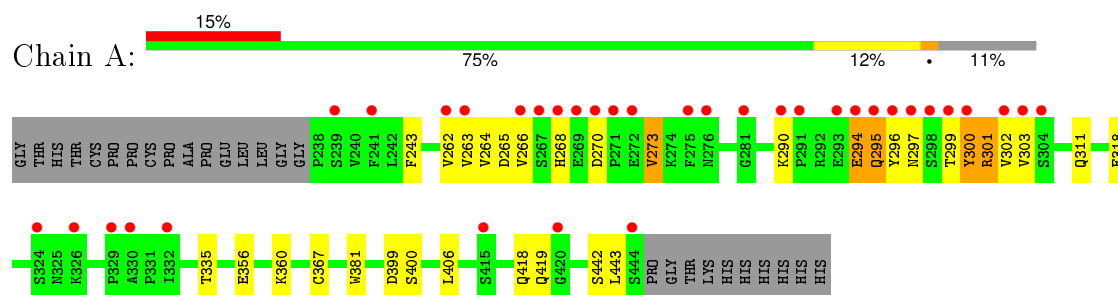
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	62	Total	O	0	0
			62	62		
5	B	104	Total	O	0	0
			104	104		

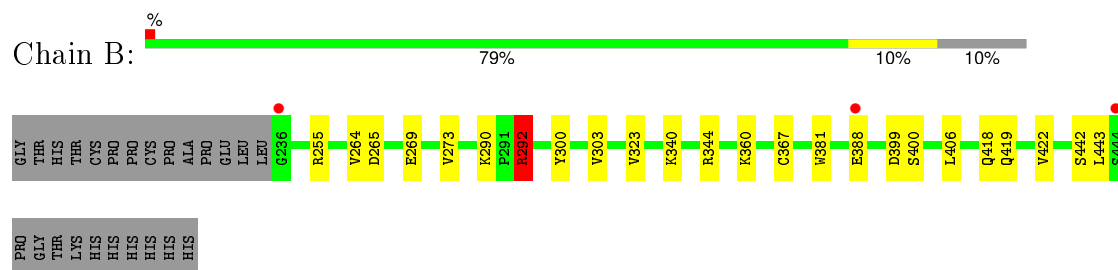
3 Residue-property plots [i](#)

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



• Molecule 1: IG GAMMA-1 CHAIN C REGION



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	153.00Å 153.00Å 111.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	66.34 – 2.30 66.25 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.0 (66.34-2.30) 99.0 (66.25-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.222 , 0.262 0.226 , 0.260	Depositor DCC
R_{free} test set	1745 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	47.0	Xtriage
Anisotropy	0.575	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 34428 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3780	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: BMA, NAG, CL, SIA, GAL, FUC, 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	0.44	0/1707	0.63	1/2324 (0.0%)
1	B	0.56	0/1758	0.71	2/2393 (0.1%)
All	All	0.50	0/3465	0.67	3/4717 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	292	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	B	292	ARG	NE-CZ-NH2	-6.00	117.30	120.30
1	A	270	ASP	CB-CG-OD1	5.03	122.83	118.30

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	1661	0	1628	53	0
1	B	1700	0	1677	15	0
2	A	130	0	110	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	B	121	0	103	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	62	0	0	2	0
5	B	104	0	0	2	0
All	All	3780	0	3518	70	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 10.

All (70) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:VAL:CG1	1:A:301:ARG:HH21	1.38	1.35
1:A:295:GLN:HB3	1:A:296:TYR:CD1	1.66	1.27
1:A:295:GLN:HB3	1:A:296:TYR:CE1	1.75	1.22
1:A:295:GLN:HG2	1:A:296:TYR:CE1	1.84	1.12
1:A:295:GLN:CB	1:A:296:TYR:CD1	2.32	1.12
1:A:262:VAL:HG11	1:A:301:ARG:HH21	1.10	1.10
1:A:263:VAL:O	1:A:301:ARG:HA	1.59	1.02
1:A:262:VAL:CG1	1:A:301:ARG:NH2	2.22	1.01
1:A:295:GLN:HG2	1:A:296:TYR:HE1	1.15	1.00
1:A:295:GLN:CB	1:A:296:TYR:CE1	2.44	0.99
1:A:295:GLN:CG	1:A:296:TYR:CE1	2.48	0.96
1:A:268:HIS:HA	1:A:300:TYR:OH	1.68	0.93
1:A:262:VAL:HG13	1:A:301:ARG:HH21	1.33	0.92
1:A:262:VAL:HG11	1:A:301:ARG:NH2	1.84	0.87
2:A:1445:NAG:H81	5:A:2059:HOH:O	1.79	0.82
1:A:318:GLU:HG2	1:A:335:THR:CG2	2.11	0.80
1:A:273:VAL:HG21	1:A:302:VAL:HG21	1.62	0.80
1:A:262:VAL:HG13	1:A:301:ARG:NH2	1.95	0.77
1:A:295:GLN:CG	1:A:296:TYR:CD1	2.71	0.71
1:A:295:GLN:CG	1:A:296:TYR:HE1	1.94	0.69
1:A:265:ASP:HA	1:A:299:THR:OG1	1.93	0.68
1:B:290:LYS:HE2	1:B:303[A]:VAL:CG2	2.24	0.68
1:A:263:VAL:O	1:A:301:ARG:CA	2.40	0.67
1:B:418:GLN:HA	1:B:443:LEU:HD22	1.80	0.64
1:A:295:GLN:CA	1:A:296:TYR:CD1	2.82	0.62
1:A:266:VAL:O	1:A:300:TYR:CD1	2.53	0.61
1:A:418:GLN:HA	1:A:443:LEU:HD22	1.81	0.61
1:A:265:ASP:HA	1:A:299:THR:HG21	1.82	0.60
1:A:295:GLN:HG2	1:A:296:TYR:CD1	2.30	0.60
1:A:266:VAL:O	1:A:300:TYR:CE1	2.56	0.59
1:B:292:ARG:HD3	5:B:2034:HOH:O	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:ASN:OD1	1:A:297:ASN:C	2.40	0.58
1:B:290:LYS:HE2	1:B:303[A]:VAL:HG21	1.83	0.58
1:A:262:VAL:HG22	1:A:303:VAL:HG13	1.84	0.58
1:A:264:VAL:HG12	1:A:301:ARG:HB3	1.87	0.56
1:A:265:ASP:HA	1:A:299:THR:CB	2.36	0.54
1:A:266:VAL:N	1:A:299:THR:HB	2.21	0.54
1:A:300:TYR:N	1:A:300:TYR:CD1	2.75	0.54
1:A:262:VAL:CG2	1:A:301:ARG:NH2	2.71	0.53
1:B:292:ARG:HG3	1:B:300:TYR:CD1	2.42	0.53
1:B:422:VAL:HG22	1:B:442:SER:OG	2.10	0.52
4:B:1452:NAG:H83	5:B:2100:HOH:O	2.09	0.52
1:A:356:GLU:HG2	5:A:2030:HOH:O	2.11	0.51
1:A:265:ASP:HA	1:A:299:THR:CG2	2.40	0.50
1:B:418:GLN:O	1:B:419:GLN:HG2	2.10	0.50
1:A:290:LYS:HD3	1:A:303:VAL:HB	1.92	0.50
1:A:418:GLN:O	1:A:419:GLN:HG2	2.12	0.50
1:A:295:GLN:HA	1:A:296:TYR:HA	1.55	0.48
1:A:295:GLN:HA	1:A:296:TYR:CD1	2.50	0.47
1:A:418:GLN:O	1:A:419:GLN:CG	2.65	0.45
1:A:265:ASP:CA	1:A:299:THR:OG1	2.64	0.45
1:B:418:GLN:O	1:B:419:GLN:CG	2.64	0.45
1:A:266:VAL:HG12	1:A:300:TYR:CE1	2.52	0.44
1:B:269[A]:GLU:H	1:B:269[A]:GLU:CD	2.19	0.44
1:B:264:VAL:O	1:B:265:ASP:HB2	2.16	0.44
1:A:295:GLN:HA	1:A:296:TYR:HD1	1.83	0.43
1:A:418:GLN:O	1:A:419:GLN:CB	2.66	0.43
1:A:406:LEU:HD12	1:A:406:LEU:C	2.39	0.43
1:A:262:VAL:CG2	1:A:301:ARG:HH21	2.31	0.43
1:A:290:LYS:HG3	1:A:303:VAL:O	2.19	0.42
1:A:243:PHE:CG	2:A:1451:NAG:H5	2.54	0.42
1:B:406:LEU:HD12	1:B:406:LEU:C	2.39	0.42
1:B:418:GLN:O	1:B:419:GLN:CB	2.66	0.42
1:A:262:VAL:CB	1:A:301:ARG:HH21	2.19	0.42
1:A:266:VAL:HB	1:A:300:TYR:O	2.20	0.41
1:B:367:CYS:HB2	1:B:381:TRP:CZ2	2.55	0.41
1:A:294:GLU:HG2	1:A:300:TYR:HD2	1.86	0.41
1:A:367:CYS:HB2	1:A:381:TRP:CZ2	2.56	0.41
1:B:273:VAL:HG13	1:B:323:VAL:HG13	2.03	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	205/233 (88%)	200 (98%)	5 (2%)	0	100	100
1	B	212/233 (91%)	208 (98%)	4 (2%)	0	100	100
All	All	417/466 (90%)	408 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	194/215 (90%)	184 (95%)	10 (5%)	29	38
1	B	199/215 (93%)	190 (96%)	9 (4%)	34	46
All	All	393/430 (91%)	374 (95%)	19 (5%)	33	42

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	273	VAL
1	A	294	GLU
1	A	295	GLN
1	A	300	TYR
1	A	301	ARG
1	A	311	GLN
1	A	360	LYS
1	A	399	ASP

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Mol	Chain	Res	Type
1	A	400	SER
1	A	442	SER
1	B	255[A]	ARG
1	B	255[B]	ARG
1	B	292	ARG
1	B	340	LYS
1	B	344	ARG
1	B	360	LYS
1	B	388	GLU
1	B	399	ASP
1	B	400	SER

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

Mol	Chain	Res	Type
1	A	421	ASN

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 ⓘ

20 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	1445	1,2	14,14,15	0.66	0	15,19,21	1.45	3 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	A	1446	2	14,14,15	0.73	0	15,19,21	1.35	3 (20%)
2	BMA	A	1447	2	11,11,12	0.54	0	14,15,17	1.65	3 (21%)
2	MAN	A	1448	2	11,11,12	0.67	0	14,15,17	0.98	0
2	NAG	A	1449	2	14,14,15	0.44	0	15,19,21	1.23	1 (6%)
2	MAN	A	1450	2	11,11,12	0.63	0	14,15,17	1.19	0
2	NAG	A	1451	2	14,14,15	0.73	0	15,19,21	0.64	0
2	GAL	A	1452	2	11,11,12	0.69	0	14,15,17	0.83	1 (7%)
2	SIA	A	1453	2	16,20,21	0.65	0	18,28,31	1.46	3 (16%)
2	FUC	A	1454	2	10,10,11	0.83	0	14,14,16	1.12	2 (14%)
4	NAG	B	1445	1,4	14,14,15	0.44	0	15,19,21	0.98	1 (6%)
4	NAG	B	1446	4	14,14,15	0.76	0	15,19,21	0.67	0
4	BMA	B	1447	4	11,11,12	0.51	0	14,15,17	1.50	4 (28%)
4	MAN	B	1448	4	11,11,12	0.60	0	14,15,17	1.16	1 (7%)
4	NAG	B	1449	4	14,14,15	0.60	0	15,19,21	1.48	3 (20%)
4	GAL	B	1450	4	11,11,12	0.49	0	14,15,17	0.83	0
4	MAN	B	1451	4	11,11,12	0.75	0	14,15,17	1.79	2 (14%)
4	NAG	B	1452	4	14,14,15	0.66	0	15,19,21	1.42	3 (20%)
4	GAL	B	1453	4	11,11,12	0.67	0	14,15,17	1.44	3 (21%)
4	FUC	B	1454	4	10,10,11	0.89	0	14,14,16	2.68	4 (28%)

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	1445	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1446	2	-	0/6/23/26	0/1/1/1
2	BMA	A	1447	2	-	0/2/19/22	0/1/1/1
2	MAN	A	1448	2	-	0/2/19/22	0/1/1/1
2	NAG	A	1449	2	-	0/6/23/26	0/1/1/1
2	MAN	A	1450	2	-	0/2/19/22	0/1/1/1
2	NAG	A	1451	2	-	0/6/23/26	0/1/1/1
2	GAL	A	1452	2	-	0/2/19/22	0/1/1/1
2	SIA	A	1453	2	-	0/14/34/38	0/1/1/1
2	FUC	A	1454	2	-	0/0/17/20	0/1/1/1
4	NAG	B	1445	1,4	-	0/6/23/26	0/1/1/1
4	NAG	B	1446	4	-	0/6/23/26	0/1/1/1
4	BMA	B	1447	4	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MAN	B	1448	4	-	0/2/19/22	0/1/1/1
4	NAG	B	1449	4	-	0/6/23/26	0/1/1/1
4	GAL	B	1450	4	-	0/2/19/22	0/1/1/1
4	MAN	B	1451	4	-	0/2/19/22	0/1/1/1
4	NAG	B	1452	4	-	0/6/23/26	0/1/1/1
4	GAL	B	1453	4	-	0/2/19/22	0/1/1/1
4	FUC	B	1454	4	-	0/0/17/20	0/1/1/1

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1454	FUC	C1-C2-C3	-6.69	101.63	109.54
4	B	1454	FUC	C1-O5-C5	-5.06	104.56	112.38
4	B	1451	MAN	O2-C2-C3	-3.27	103.54	110.12
4	B	1449	NAG	O7-C7-C8	-2.83	116.86	122.06
4	B	1452	NAG	O7-C7-C8	-2.73	117.05	122.06
2	A	1453	SIA	C7-C6-C5	-2.70	110.23	114.32
2	A	1446	NAG	O7-C7-C8	-2.57	117.35	122.06
2	A	1445	NAG	O6-C6-C5	-2.52	103.01	111.33
4	B	1454	FUC	O4-C4-C5	-2.45	104.09	109.84
4	B	1447	BMA	O6-C6-C5	-2.33	103.64	111.33
2	A	1447	BMA	O3-C3-C4	-2.24	105.29	110.34
4	B	1447	BMA	O3-C3-C4	-2.06	105.69	110.34
4	B	1445	NAG	C3-C4-C5	-2.02	106.67	110.20
2	A	1445	NAG	C8-C7-N2	2.00	119.94	116.11
2	A	1452	GAL	C1-O5-C5	2.08	114.89	112.25
4	B	1452	NAG	C1-O5-C5	2.10	114.91	112.25
4	B	1453	GAL	C3-C4-C5	2.11	113.87	110.20
4	B	1453	GAL	O2-C2-C1	2.21	113.64	109.21
2	A	1446	NAG	O5-C5-C6	2.22	112.16	107.35
4	B	1447	BMA	C1-O5-C5	2.23	115.07	112.25
4	B	1453	GAL	C1-O5-C5	2.24	115.09	112.25
2	A	1454	FUC	O5-C5-C6	2.25	109.85	106.13
2	A	1449	NAG	C8-C7-N2	2.36	120.62	116.11
2	A	1454	FUC	C1-C2-C3	2.45	112.44	109.54
2	A	1445	NAG	C1-O5-C5	2.61	115.56	112.25
2	A	1446	NAG	C4-C3-C2	2.67	115.37	111.23
2	A	1447	BMA	C3-C4-C5	2.69	114.89	110.20
4	B	1449	NAG	C2-N2-C7	2.76	126.59	123.04
4	B	1447	BMA	C1-C2-C3	2.79	112.84	109.54
4	B	1448	MAN	C1-O5-C5	2.85	115.86	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1449	NAG	C8-C7-N2	3.13	122.10	116.11
2	A	1453	SIA	O6-C6-C5	3.33	113.94	108.48
2	A	1453	SIA	C3-C4-C5	3.35	115.21	111.47
4	B	1452	NAG	C8-C7-N2	3.42	122.66	116.11
2	A	1447	BMA	C1-O5-C5	3.45	116.62	112.25
4	B	1454	FUC	O5-C5-C6	4.14	112.97	106.13
4	B	1451	MAN	C1-C2-C3	4.36	114.70	109.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1445	NAG	1	0
2	A	1451	NAG	1	0
4	B	1452	NAG	1	0

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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 ⓘ

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	207/233 (88%)	0.93	35 (16%) 2 3	40, 71, 141, 165	0
1	B	209/233 (89%)	0.33	3 (1%) 78 83	31, 47, 108, 131	1 (0%)
All	All	416/466 (89%)	0.63	38 (9%) 11 17	31, 59, 128, 165	1 (0%)

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	302	VAL	11.2
1	A	300	TYR	8.2
1	A	295	GLN	6.6
1	A	266	VAL	6.4
1	A	296	TYR	6.3
1	A	269	GLU	5.8
1	A	298	SER	4.5
1	A	299	THR	4.3
1	A	267	SER	4.0
1	A	303	VAL	3.9
1	A	271	PRO	3.8
1	A	241	PHE	3.8
1	A	263	VAL	3.7
1	A	291	PRO	3.6
1	A	272	GLU	3.5
1	A	281	GLY	3.5
1	A	297	ASN	3.5
1	A	290	LYS	3.4
1	A	420	GLY	3.2
1	A	326	LYS	3.2
1	A	329	PRO	3.2
1	B	236	GLY	3.1
1	A	444	SER	3.0
1	A	262	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	276	ASN	2.9
1	A	239	SER	2.9
1	A	268	HIS	2.8
1	A	293	GLU	2.8
1	A	294	GLU	2.7
1	A	330	ALA	2.6
1	A	270	ASP	2.6
1	A	324	SER	2.6
1	A	415	SER	2.6
1	A	275	PHE	2.4
1	A	332	ILE	2.4
1	A	304	SER	2.3
1	B	444	SER	2.1
1	B	388	GLU	2.1

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

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

6.3 Carbohydrates [i](#)

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	GAL	A	1452	11/12	0.83	0.12	-0.83	58,63,72,89	0
4	NAG	B	1445	14/15	0.97	0.12	-1.49	32,37,40,42	0
4	NAG	B	1452	14/15	0.94	0.12	-1.69	44,46,54,59	0
4	GAL	B	1453	11/12	0.96	0.13	-2.51	43,47,55,61	0
4	GAL	B	1450	11/12	0.60	0.31	-	111,121,130,134	0
2	BMA	A	1447	11/12	0.87	0.12	-	64,69,73,74	0
2	NAG	A	1451	14/15	0.83	0.20	-	64,70,75,76	0
2	NAG	A	1449	14/15	0.69	0.27	-	116,126,138,140	0
4	FUC	B	1454	10/11	0.93	0.13	-	42,45,48,51	0
2	NAG	A	1445	14/15	0.89	0.18	-	84,91,97,98	0
2	MAN	A	1448	11/12	0.91	0.14	-	80,89,93,108	0
2	NAG	A	1446	14/15	0.94	0.17	-	71,81,88,88	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MAN	A	1450	11/12	0.87	0.17	-	74,77,84,94	0
4	NAG	B	1449	14/15	0.70	0.19	-	98,102,117,129	0
2	SIA	A	1453	20/21	0.72	0.38	-	122,131,136,136	0
4	MAN	B	1451	11/12	0.96	0.13	-	42,49,55,59	0
4	NAG	B	1446	14/15	0.97	0.10	-	27,36,44,45	0
4	BMA	B	1447	11/12	0.94	0.12	-	41,44,52,55	0
2	FUC	A	1454	10/11	0.91	0.21	-	94,103,106,108	0
4	MAN	B	1448	11/12	0.92	0.13	-	64,73,81,83	0

6.4 Ligands [i](#)

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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	CL	B	1455	1/1	1.00	0.18	-	38,38,38,38	1
3	CL	A	1455	1/1	0.99	0.09	-	44,44,44,44	1

6.5 Other polymers [i](#)

There are no such residues in this entry.