



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 06:56 am GMT

PDB ID : 3D6G
Title : Fc fragment of IgG1 (Herceptin) with protein-A mimetic peptide dendrimer ligand.
Authors : Bujacz, A.D.; Redzynia, I.; Bujacz, G.D.; Dinon, F.; Pengo, P.; Fassina, G.
Deposited on : 2008-05-19
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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

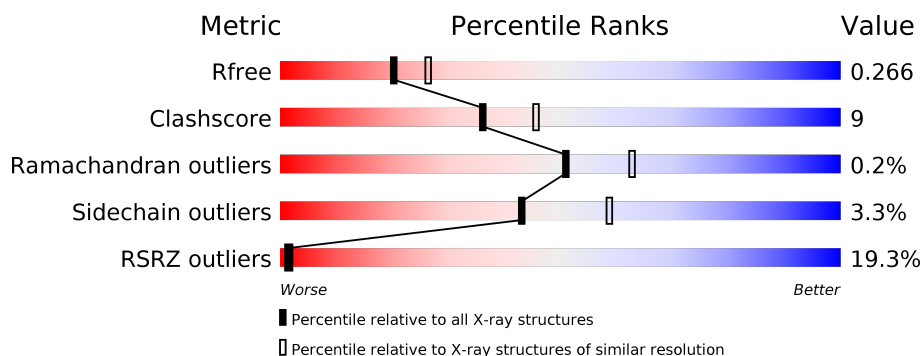
1 Overall quality at a glance

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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (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	210	<div> <div>12%</div> <div>90%</div> <div>10%</div> </div>
1	B	210	<div> <div>27%</div> <div>79%</div> <div>19%</div> <div>.</div> </div>

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	X12	B	446	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4002 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.

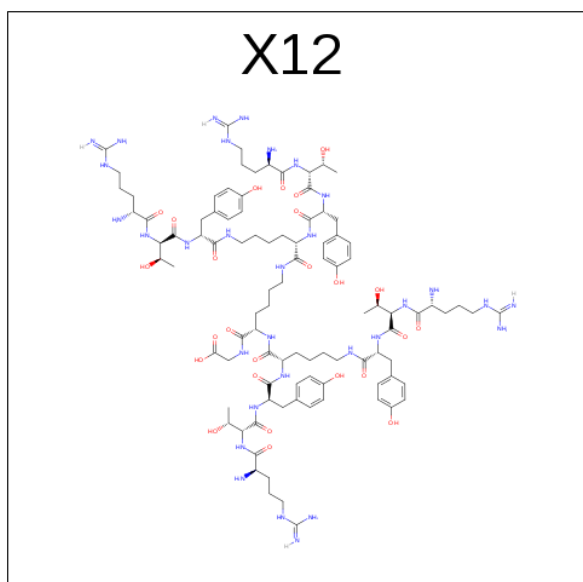
- Molecule 1 is a protein called Ig gamma-1 chain C region.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	210	Total	C	N	O	S	0	4	0
			1702	1084	288	324	6			
1	B	209	Total	C	N	O	S	0	2	0
			1682	1070	285	321	6			

- Molecule 2 is a polymer of unknown type called SUGAR (8-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	8	Total	C	N	O	0	0
			99	56	4	39		
2	B	8	Total	C	N	O	0	0
			99	56	4	39		

- Molecule 3 is 2-[[[(2S)-2,6-BIS[[[(2S)-2,6-BIS[[[(2R)-2-[[[(2R,3R)-2-[[[(2R)-2-AMINO-5-CARBA MIMIDAMIDO-PENTANOYL]AMINO]-3-HYDROXY-BUTANOYL]AMINO]-3-(4-HYDROXYPHENYL)PROPANOYL]AMINO]HEXANOYL]AMINO]HEXANOYL]AMINO]ETHANOIC ACID (three-letter code: X12) (formula: C₉₆H₁₅₃N₃₁O₂₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			152	96	31	25		

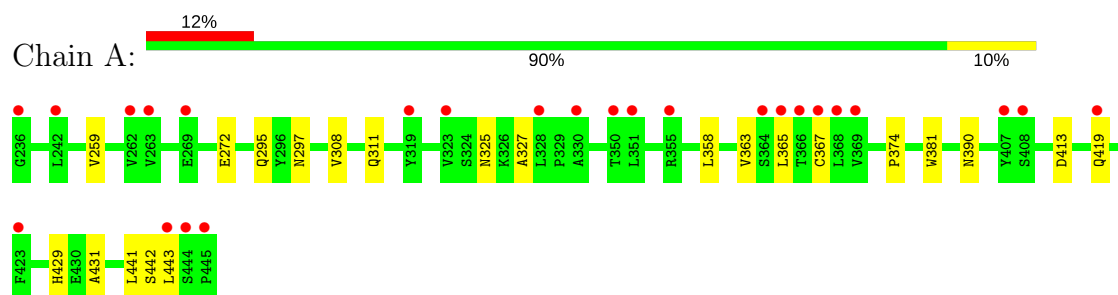
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	145	Total	O	0	0
			145	145		
4	B	123	Total	O	0	0
			123	123		

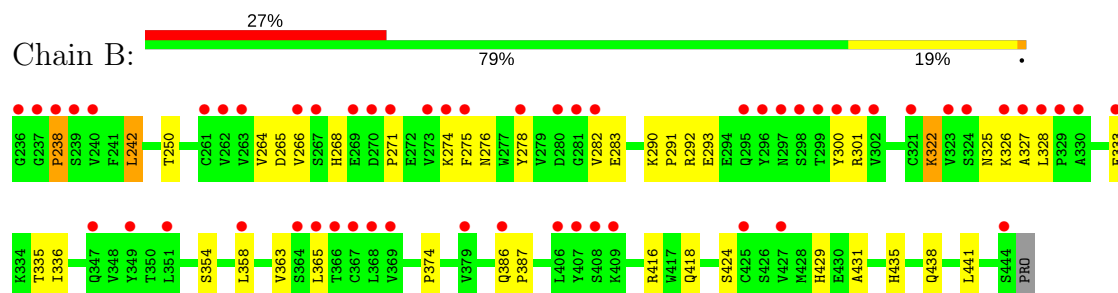
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.78Å 79.16Å 139.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.30 30.11 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.3 (40.00-2.30) 99.3 (30.11-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.188 , 0.257 0.199 , 0.266	Depositor DCC
R_{free} test set	806 reflections (3.33%)	DCC
Wilson B-factor (Å ²)	53.1	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 62.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4002	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.78% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, X12, BMA, NAG, FUC

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	0.76	0/1762	0.77	0/2400
1	B	0.76	0/1734	0.74	0/2361
All	All	0.76	0/3496	0.75	0/4761

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1702	0	1674	12	0
1	B	1682	0	1654	43	0
2	A	99	0	85	2	0
2	B	99	0	85	2	0
3	B	152	0	146	13	0
4	A	145	0	0	0	0
4	B	123	0	0	4	0
All	All	4002	0	3644	67	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 9.

All (67) 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)
3:B:446:X12:HNBS	3:B:446:X12:HBI	1.29	0.96
1:B:276:ASN:HA	4:B:532:HOH:O	1.65	0.95
1:B:282:VAL:HG12	1:B:283:GLU:N	1.91	0.83
3:B:446:X12:HBI	3:B:446:X12:NBS	1.93	0.83
1:B:429:HIS:HD2	1:B:431:ALA:H	1.32	0.77
1:B:274:LYS:HG2	1:B:275:PHE:N	2.01	0.76
1:B:278:TYR:HD1	1:B:282:VAL:O	1.73	0.72
1:B:282:VAL:HG12	1:B:283:GLU:O	1.90	0.72
1:B:274:LYS:HG2	1:B:275:PHE:H	1.55	0.71
1:B:278:TYR:CD1	1:B:282:VAL:O	2.44	0.70
1:B:242:LEU:HD13	1:B:336:ILE:HB	1.72	0.70
1:A:297:ASN:HD22	2:A:1:NAG:H83	1.58	0.69
1:B:429:HIS:CD2	1:B:431:ALA:H	2.12	0.65
1:B:268:HIS:CD2	1:B:300:TYR:CE1	2.86	0.64
1:B:322:LYS:HG2	1:B:333:GLU:HG2	1.79	0.64
1:B:264:VAL:HG13	1:B:265:ASP:H	1.63	0.64
1:B:268:HIS:CD2	1:B:300:TYR:HE1	2.18	0.61
1:B:278:TYR:HB3	1:B:282:VAL:O	1.99	0.61
1:B:264:VAL:HG13	1:B:265:ASP:N	2.17	0.60
1:B:358:LEU:HD23	1:B:363:VAL:HG11	1.84	0.60
1:B:282:VAL:CG1	1:B:283:GLU:N	2.61	0.59
1:A:374:PRO:O	1:A:429:HIS:HE1	1.86	0.59
1:B:250:THR:O	3:B:446:X12:NDU	2.36	0.58
1:B:325:ASN:HB3	1:B:328:LEU:HG	1.88	0.56
1:B:278:TYR:CB	1:B:282:VAL:O	2.56	0.54
1:B:282:VAL:HG12	1:B:283:GLU:H	1.70	0.53
1:A:272[A]:GLU:HA	1:A:272[A]:GLU:OE1	2.08	0.53
1:B:264:VAL:CG1	1:B:265:ASP:N	2.71	0.53
1:A:429:HIS:CD2	1:A:431:ALA:H	2.27	0.52
3:B:446:X12:HCQ	3:B:446:X12:HB5	1.92	0.52
1:A:295:GLN:HE21	2:A:1:NAG:H62	1.75	0.51
1:A:429:HIS:HD2	1:A:431:ALA:H	1.57	0.51
1:B:274:LYS:CG	1:B:275:PHE:N	2.74	0.50
3:B:446:X12:HCJ	3:B:446:X12:HC6	1.93	0.50
1:A:259:VAL:HG23	1:A:308:VAL:HG11	1.94	0.50
1:A:367:CYS:HB2	1:A:381:TRP:CZ2	2.48	0.49
1:B:276:ASN:HB2	1:B:322:LYS:HB2	1.95	0.48
1:B:325:ASN:HD21	1:B:327:ALA:HB3	1.79	0.48
1:A:311[B]:GLN:H	1:A:311[B]:GLN:NE2	2.11	0.48
3:B:446:X12:HAW	3:B:446:X12:NBC	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:LYS:CG	1:B:275:PHE:H	2.26	0.46
3:B:446:X12:HAT	3:B:446:X12:HAQA	1.48	0.45
3:B:446:X12:HNAG	3:B:446:X12:CC1	2.28	0.45
1:B:435:HIS:HE1	3:B:446:X12:NDT	2.14	0.45
1:B:424:SER:HB3	1:B:438:GLN:HE21	1.82	0.45
1:B:266:VAL:HB	1:B:300:TYR:HB2	2.00	0.44
1:B:325:ASN:ND2	1:B:327:ALA:H	2.15	0.44
1:A:358:LEU:HD23	1:A:363:VAL:HG11	1.98	0.44
1:B:282:VAL:CG1	1:B:283:GLU:H	2.29	0.43
1:B:325:ASN:ND2	1:B:327:ALA:HB3	2.34	0.43
2:B:3:BMA:O4	4:B:512:HOH:O	2.21	0.42
1:B:365:LEU:HB3	1:B:441:LEU:HD23	2.01	0.42
1:B:276:ASN:HB2	1:B:322:LYS:CB	2.48	0.42
1:B:335:THR:HG22	1:B:336:ILE:N	2.34	0.42
1:B:416:ARG:HD3	4:B:509:HOH:O	2.18	0.42
1:B:238:PRO:HB3	1:B:264:VAL:O	2.20	0.41
1:B:271:PRO:HG2	4:B:560:HOH:O	2.20	0.41
1:B:374:PRO:O	1:B:429:HIS:HE1	2.02	0.41
1:B:435:HIS:HE1	3:B:446:X12:CDS	2.33	0.41
1:B:386:GLN:HA	1:B:387:PRO:HD3	1.96	0.41
1:B:290:LYS:HA	1:B:291:PRO:HD3	1.79	0.41
3:B:446:X12:HCU	3:B:446:X12:HDH	2.01	0.41
1:A:325:ASN:OD1	1:A:327:ALA:HB3	2.21	0.40
3:B:446:X12:HBP	3:B:446:X12:HNBA	1.86	0.40
2:B:4:MAN:H4	2:B:5:NAG:H83	2.03	0.40
1:A:365:LEU:HB3	1:A:441:LEU:HD23	2.02	0.40
3:B:446:X12:CAW	3:B:446:X12:NBC	2.84	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	212/210 (101%)	210 (99%)	2 (1%)	0	100	100
1	B	209/210 (100%)	202 (97%)	6 (3%)	1 (0%)	32	39
All	All	421/420 (100%)	412 (98%)	8 (2%)	1 (0%)	51	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	238	PRO

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	199/195 (102%)	194 (98%)	5 (2%)	53	70
1	B	196/195 (100%)	187 (95%)	9 (5%)	31	42
All	All	395/390 (101%)	381 (96%)	14 (4%)	43	56

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	390	ASN
1	A	413	ASP
1	A	419	GLN
1	A	442	SER
1	A	443	LEU
1	B	242	LEU
1	B	292	ARG
1	B	293	GLU
1	B	301[A]	ARG
1	B	301[B]	ARG
1	B	322	LYS
1	B	326	LYS
1	B	354	SER
1	B	418	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	295	GLN
1	A	361	ASN
1	A	386	GLN
1	A	419	GLN
1	A	429	HIS
1	A	438	GLN
1	B	268	HIS
1	B	325	ASN
1	B	389	ASN
1	B	390	ASN
1	B	418	GLN
1	B	419	GLN
1	B	429	HIS
1	B	435	HIS
1	B	438	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 ⓘ

16 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	1	1,2	14,14,15	1.31	2 (14%)	15,19,21	1.60	3 (20%)
2	NAG	A	2	2	14,14,15	1.04	1 (7%)	15,19,21	1.97	6 (40%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BMA	A	3	2	11,11,12	1.14	2 (18%)	13,15,17	1.62	2 (15%)
2	MAN	A	4	2	11,11,12	0.98	1 (9%)	13,15,17	2.04	4 (30%)
2	NAG	A	5	2	14,14,15	0.95	1 (7%)	15,19,21	1.23	2 (13%)
2	MAN	A	6	2	11,11,12	1.36	2 (18%)	13,15,17	1.26	2 (15%)
2	NAG	A	7	2	14,14,15	0.79	1 (7%)	15,19,21	1.56	3 (20%)
2	FUC	A	8	2	9,10,11	0.43	0	13,14,16	0.98	0
2	NAG	B	1	1,2	14,14,15	1.25	2 (14%)	15,19,21	1.46	2 (13%)
2	NAG	B	2	2	14,14,15	0.96	0	15,19,21	1.49	2 (13%)
2	BMA	B	3	2	11,11,12	0.86	1 (9%)	13,15,17	1.53	3 (23%)
2	MAN	B	4	2	11,11,12	1.10	1 (9%)	13,15,17	1.75	2 (15%)
2	NAG	B	5	2	14,14,15	1.09	1 (7%)	15,19,21	1.42	2 (13%)
2	MAN	B	6	2	11,11,12	1.11	1 (9%)	13,15,17	1.67	1 (7%)
2	NAG	B	7	2	14,14,15	0.79	0	15,19,21	1.08	1 (6%)
2	FUC	B	8	2	9,10,11	0.60	0	13,14,16	1.47	1 (7%)

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	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	2	2	-	0/6/23/26	0/1/1/1
2	BMA	A	3	2	-	0/2/19/22	0/1/1/1
2	MAN	A	4	2	-	0/2/19/22	0/1/1/1
2	NAG	A	5	2	-	0/6/23/26	0/1/1/1
2	MAN	A	6	2	-	0/2/19/22	0/1/1/1
2	NAG	A	7	2	-	0/6/23/26	0/1/1/1
2	FUC	A	8	2	-	0/0/17/20	0/1/1/1
2	NAG	B	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	BMA	B	3	2	-	0/2/19/22	0/1/1/1
2	MAN	B	4	2	-	0/2/19/22	0/1/1/1
2	NAG	B	5	2	-	0/6/23/26	0/1/1/1
2	MAN	B	6	2	-	0/2/19/22	0/1/1/1
2	NAG	B	7	2	-	0/6/23/26	0/1/1/1
2	FUC	B	8	2	-	0/0/17/20	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2	NAG	O5-C1	-2.45	1.39	1.43
2	B	3	BMA	C2-C3	2.02	1.55	1.52
2	B	5	NAG	C4-C5	2.08	1.57	1.53
2	A	3	BMA	O5-C5	2.12	1.47	1.43
2	A	6	MAN	O5-C5	2.15	1.48	1.43
2	A	7	NAG	C8-C7	2.17	1.55	1.50
2	A	5	NAG	C8-C7	2.37	1.55	1.50
2	B	4	MAN	O5-C1	2.38	1.47	1.43
2	B	1	NAG	O5-C1	2.38	1.47	1.43
2	B	1	NAG	C1-C2	2.43	1.55	1.52
2	A	4	MAN	O5-C5	2.45	1.48	1.43
2	A	3	BMA	C2-C3	2.47	1.55	1.52
2	A	1	NAG	C1-C2	2.65	1.56	1.52
2	B	6	MAN	O5-C1	2.68	1.48	1.43
2	A	1	NAG	O5-C5	2.79	1.49	1.43
2	A	6	MAN	C2-C3	3.51	1.57	1.52

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	NAG	O6-C6-C5	-4.50	96.19	111.34
2	B	4	MAN	O2-C2-C3	-4.40	101.53	110.17
2	A	2	NAG	C3-C4-C5	-4.36	102.53	110.22
2	A	7	NAG	O5-C1-C2	-4.00	105.91	111.47
2	A	4	MAN	O2-C2-C3	-3.80	102.71	110.17
2	B	5	NAG	C1-C2-N2	-3.71	104.15	110.49
2	A	3	BMA	O6-C6-C5	-3.71	98.85	111.34
2	B	8	FUC	O5-C1-C2	-3.70	104.99	110.79
2	A	2	NAG	O5-C1-C2	-3.22	107.00	111.47
2	A	1	NAG	C6-C5-C4	-3.18	105.56	113.00
2	A	3	BMA	O3-C3-C4	-3.07	103.67	110.36
2	A	4	MAN	C2-C3-C4	-2.81	105.97	110.88
2	B	1	NAG	C6-C5-C4	-2.65	106.80	113.00
2	A	2	NAG	O4-C4-C5	-2.54	102.90	109.28
2	A	2	NAG	C8-C7-N2	-2.50	111.59	116.11
2	B	3	BMA	O6-C6-C5	-2.48	103.00	111.34
2	B	3	BMA	C6-C5-C4	-2.39	107.42	113.00
2	B	2	NAG	O4-C4-C5	-2.33	103.40	109.28
2	B	7	NAG	O5-C1-C2	-2.31	108.25	111.47
2	A	5	NAG	C1-O5-C5	-2.10	109.27	112.17
2	A	7	NAG	C1-O5-C5	-2.04	109.36	112.17
2	A	1	NAG	C3-C4-C5	-2.03	106.65	110.22
2	A	6	MAN	O2-C2-C1	-2.00	105.11	109.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	5	NAG	C2-N2-C7	2.09	125.99	122.94
2	B	5	NAG	O5-C1-C2	2.25	114.61	111.47
2	A	2	NAG	O7-C7-N2	2.42	126.58	121.92
2	A	4	MAN	C1-C2-C3	2.55	112.88	109.65
2	A	7	NAG	C6-C5-C4	2.67	119.24	113.00
2	A	2	NAG	C1-O5-C5	2.69	115.87	112.17
2	A	6	MAN	C1-O5-C5	2.87	116.12	112.17
2	B	3	BMA	C1-C2-C3	3.07	113.54	109.65
2	B	4	MAN	C1-C2-C3	3.59	114.20	109.65
2	B	2	NAG	C1-O5-C5	3.62	117.15	112.17
2	B	1	NAG	C1-O5-C5	3.72	117.30	112.17
2	A	4	MAN	C1-O5-C5	4.00	117.69	112.17
2	B	6	MAN	C1-O5-C5	5.52	119.77	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	NAG	2	0
2	B	3	BMA	1	0
2	B	4	MAN	1	0
2	B	5	NAG	1	0

5.6 Ligand geometry

1 ligand is 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	X12	B	446	-	152,155,155	0.96	3 (1%)	185,205,205	1.13	11 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	X12	B	446	-	-	0/191/193/193	0/4/4/4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	446	X12	CB2-NBH	2.45	1.38	1.33
3	B	446	X12	CBG-NBH	4.93	1.57	1.46
3	B	446	X12	CB2-NBK	5.20	1.53	1.32

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	446	X12	NBJ-CB2-NBH	-3.05	112.11	119.30
3	B	446	X12	CBQ-CBH-NBG	-2.01	105.67	111.20
3	B	446	X12	CBI-CBH-NBG	2.02	115.06	110.80
3	B	446	X12	CB5-NBW-CCN	2.78	127.76	121.66
3	B	446	X12	CCU-NCT-CDK	2.88	127.97	121.66
3	B	446	X12	CBT-CBX-NBG	3.12	123.27	116.53
3	B	446	X12	CAP-NAO-CBE	3.24	128.76	121.66
3	B	446	X12	CAT-NAU-C	3.25	128.66	122.59
3	B	446	X12	CAW-NAV-CB1	3.36	130.46	121.87
3	B	446	X12	CDG-NDF-CDV	3.55	130.94	121.87
3	B	446	X12	CBH-NBG-CBX	6.08	134.99	121.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	446	X12	13	0

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	210/210 (100%)	0.67	25 (11%) 5 7	65, 71, 79, 86	0
1	B	209/210 (99%)	1.32	56 (26%) 1 1	65, 72, 78, 84	0
All	All	419/420 (99%)	0.99	81 (19%) 1 2	65, 72, 79, 86	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	296	TYR	11.5
1	A	444	SER	7.0
1	B	300	TYR	6.3
1	B	366	THR	6.1
1	A	445	PRO	6.0
1	B	330	ALA	6.0
1	B	368	LEU	5.7
1	B	299	THR	5.3
1	B	240	VAL	4.8
1	B	351	LEU	4.8
1	B	323	VAL	4.7
1	B	297	ASN	4.7
1	B	281	GLY	4.7
1	B	295	GLN	4.5
1	B	367	CYS	4.3
1	B	278	TYR	4.3
1	B	408	SER	4.3
1	B	407	TYR	4.2
1	A	366	THR	4.0
1	A	368	LEU	4.0
1	B	406	LEU	3.9
1	B	236	GLY	3.9
1	B	444	SER	3.8
1	B	326	LYS	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	267	SER	3.7
1	B	270	ASP	3.7
1	A	407	TYR	3.7
1	B	369	VAL	3.7
1	B	379	VAL	3.6
1	B	327	ALA	3.6
1	B	238	PRO	3.6
1	A	355[A]	ARG	3.6
1	B	237	GLY	3.5
1	A	365	LEU	3.4
1	B	324	SER	3.4
1	B	302	VAL	3.4
1	A	269	GLU	3.3
1	B	365	LEU	3.2
1	B	301[A]	ARG	3.1
1	B	427	VAL	3.1
1	B	266	VAL	3.0
1	B	358	LEU	3.0
1	B	329	PRO	3.0
1	B	321	CYS	2.9
1	A	364	SER	2.9
1	A	328	LEU	2.9
1	B	364	SER	2.9
1	A	369	VAL	2.8
1	B	273	VAL	2.8
1	B	386	GLN	2.7
1	B	280	ASP	2.7
1	B	275	PHE	2.7
1	B	282	VAL	2.6
1	B	262	VAL	2.6
1	B	239	SER	2.6
1	B	409	LYS	2.6
1	B	333	GLU	2.5
1	A	330	ALA	2.5
1	B	271	PRO	2.5
1	A	419	GLN	2.5
1	B	269	GLU	2.4
1	A	236	GLY	2.4
1	A	263	VAL	2.4
1	A	367	CYS	2.4
1	B	298	SER	2.4
1	B	263	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	408	SER	2.3
1	B	261	CYS	2.3
1	A	262	VAL	2.3
1	B	349	TYR	2.3
1	A	443	LEU	2.2
1	A	423	PHE	2.2
1	A	351	LEU	2.2
1	A	242	LEU	2.1
1	A	323	VAL	2.1
1	A	319	TYR	2.1
1	B	274	LYS	2.1
1	B	328	LEU	2.1
1	B	425	CYS	2.1
1	B	347[A]	GLN	2.1
1	A	350	THR	2.0

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	NAG	A	7	14/15	0.79	0.17	-0.26	65,69,73,74	0
2	NAG	B	7	14/15	0.81	0.20	-0.41	64,66,68,70	0
2	MAN	A	4	11/12	0.83	0.18	-	64,68,72,77	0
2	NAG	B	2	14/15	0.45	0.35	-	62,68,71,71	0
2	MAN	A	6	11/12	0.89	0.13	-	66,69,70,72	0
2	MAN	B	6	11/12	0.85	0.19	-	65,67,69,70	0
2	BMA	B	3	11/12	0.92	0.28	-	62,65,66,68	0
2	NAG	A	5	14/15	0.68	0.24	-	82,86,89,90	0
2	FUC	A	8	10/11	0.80	0.22	-	71,75,75,76	0
2	NAG	B	1	14/15	0.59	0.48	-	59,65,66,68	0
2	BMA	A	3	11/12	0.90	0.13	-	64,68,71,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	FUC	B	8	10/11	0.85	0.54	-	69,70,71,71	0
2	NAG	A	2	14/15	0.95	0.13	-	67,69,70,73	0
2	MAN	B	4	11/12	0.88	0.13	-	61,64,67,72	0
2	NAG	B	5	14/15	0.67	0.37	-	76,78,79,79	0
2	NAG	A	1	14/15	0.92	0.10	-	66,69,74,79	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	X12	B	446	152/152	0.26	0.70	6.83	51,79,83,83	152

6.5 Other polymers [i](#)

There are no such residues in this entry.