



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 18, 2016 – 10:45 AM EDT

PDB ID : 5KJR
Title : Crystal structure of the ADCC-potent antibody N60-i3 Fab in complex with HIV-1 Clade A/E gp120 W69A/S115W mutant and M48U1.
Authors : Tolbert, W.D.; Pazgier, M.
Deposited on : 2016-06-20
Resolution : 2.98 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
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) : rb-20027939

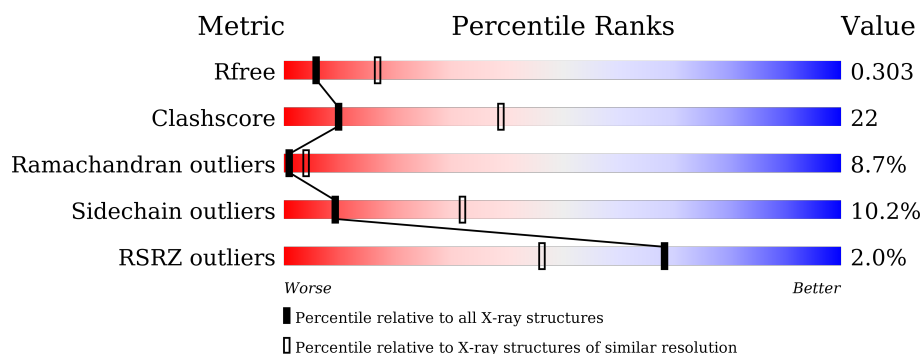
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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	1992 (3.00-2.96)
Clashscore	102246	2349 (3.00-2.96)
Ramachandran outliers	100387	2274 (3.00-2.96)
Sidechain outliers	100360	2277 (3.00-2.96)
RSRZ outliers	91569	2007 (3.00-2.96)

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	G	353	<div> <div>4%</div> <div> <div>46%</div> <div>39%</div> <div>8%</div> <div>6%</div> </div> </div>
2	N	28	<div> <div>86%</div> <div>11%</div> </div>
3	H	229	<div> <div>45%</div> <div>38%</div> <div>11%</div> <div>5%</div> </div>
4	L	221	<div> <div>%</div> <div>48%</div> <div>40%</div> <div>7%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	NAG	G	501	-	-	-	X
5	NAG	G	510	-	-	-	X
6	MPD	L	301	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6203 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 clade A/E 93TH057 HIV-1 gp120 core.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	332	Total	C	N	O	S	0	0	0
			2607	1637	452	496	22			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	69	ALA	TRP	engineered mutation	UNP A0A0M3KKW9
G	115	TRP	SER	engineered mutation	UNP A0A0M3KKW9
G	375	SER	HIS	engineered mutation	UNP A0A0M3KKW9

- Molecule 2 is a protein called M48U1 CD4 MIMETIC PEPTIDE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	N	28	Total	C	N	O	S	0	0	1
			209	133	38	32	6			

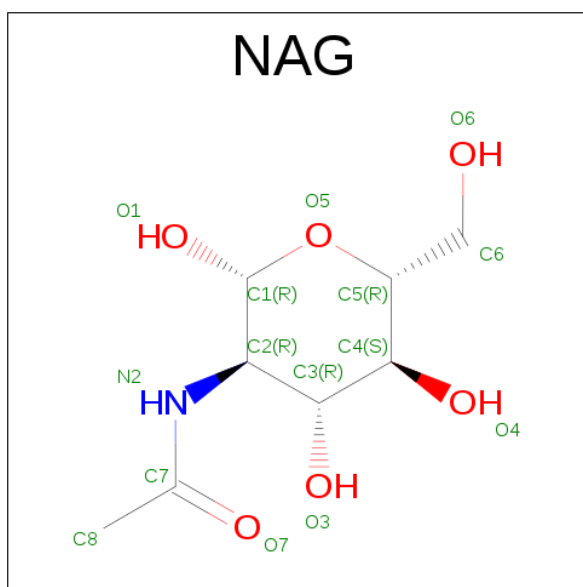
- Molecule 3 is a protein called N60-I3 FAB HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	218	Total	C	N	O	S	0	1	0
			1651	1050	277	320	4			

- Molecule 4 is a protein called N60-I3 FAB LIGHT CHAIN.

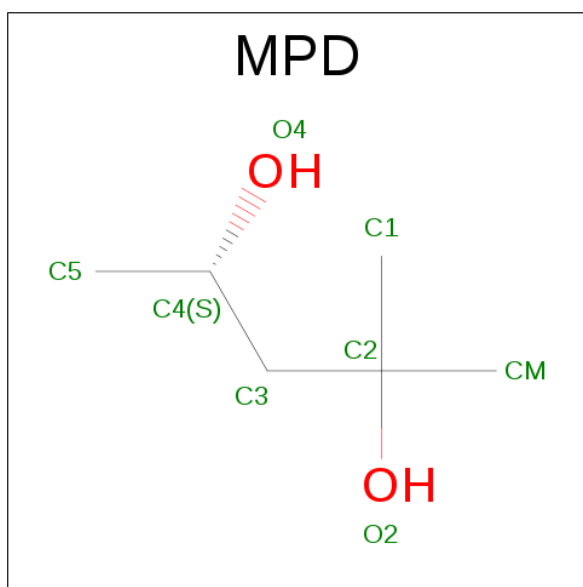
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	212	Total	C	N	O	S	0	0	0
			1579	986	261	327	5			

- Molecule 5 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	G	1	Total	C	N	O	0	0
			14	8	1	5		
5	G	1	Total	C	N	O	0	0
			14	8	1	5		
5	G	1	Total	C	N	O	0	0
			14	8	1	5		
5	G	1	Total	C	N	O	0	0
			14	8	1	5		
5	G	1	Total	C	N	O	0	0
			14	8	1	5		
5	G	1	Total	C	N	O	0	0
			14	8	1	5		
5	G	1	Total	C	N	O	0	0
			14	8	1	5		
5	G	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).

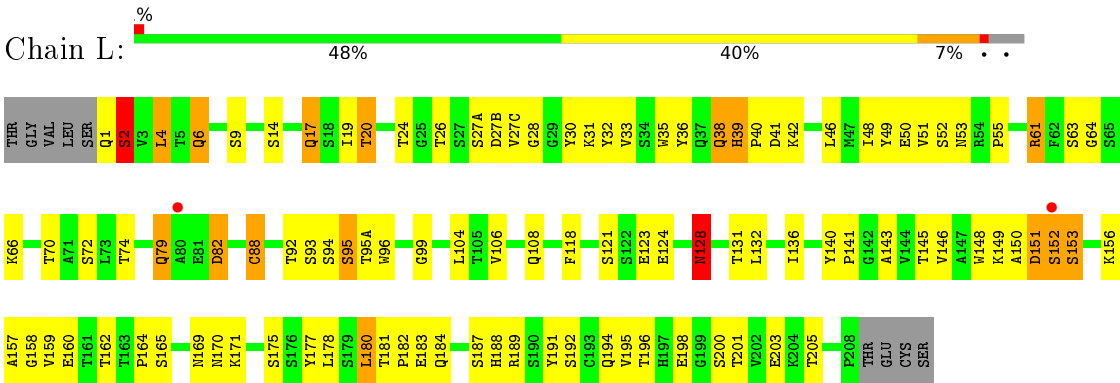


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	L	1	Total	C	O	0	0
			8	6	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	G	4	Total	O	0	0
			4	4		
7	H	3	Total	O	0	0
			3	3		
7	L	2	Total	O	0	0
			2	2		

● Molecule 4: N60-I3 FAB LIGHT CHAIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	95.19Å 101.94Å 108.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.98 42.84 – 2.98	Depositor EDS
% Data completeness (in resolution range)	87.6 (50.00-2.98) 87.7 (42.84-2.98)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.05 (at 2.95Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.249 , 0.306 0.248 , 0.303	Depositor DCC
R_{free} test set	1007 reflections (5.45%)	DCC
Wilson B-factor (Å ²)	52.3	Xtriage
Anisotropy	0.579	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 34.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	6203	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.84% 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: MPD, NAG, MPT, NH2, U2X, DPR

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	G	0.69	2/2661 (0.1%)	0.93	4/3611 (0.1%)
2	N	0.50	0/176	0.82	0/231
3	H	0.72	0/1694	0.96	4/2316 (0.2%)
4	L	0.69	1/1618 (0.1%)	0.87	1/2208 (0.0%)
All	All	0.69	3/6149 (0.0%)	0.92	9/8366 (0.1%)

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
1	G	0	2
2	N	0	2
All	All	0	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	87	GLU	CG-CD	8.85	1.65	1.51
1	G	87	GLU	CD-OE2	8.32	1.34	1.25
4	L	9	SER	C-N	5.05	1.45	1.34

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	228	CYS	CA-CB-SG	-12.55	91.42	114.00
3	H	196	CYS	CA-CB-SG	-7.26	100.93	114.00
1	G	74	CYS	CA-CB-SG	-6.47	102.34	114.00
3	H	100(C)	TYR	N-CA-C	5.87	126.85	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	461	ASN	N-CA-C	5.85	126.80	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	G	409	GLY	Peptide
1	G	462	ASN	Peptide
2	N	22	THR	Mainchain
2	N	23	U2X	Mainchain

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	G	2607	0	2537	120	0
2	N	209	0	212	1	0
3	H	1651	0	1627	96	0
4	L	1579	0	1524	70	0
5	G	140	0	130	0	0
6	L	8	0	14	0	0
7	G	4	0	0	0	0
7	H	3	0	0	0	0
7	L	2	0	0	0	0
All	All	6203	0	6044	272	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 22.

The worst 5 of 272 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:G:107:ASP:OD1	3:H:99:ARG:NH1	1.92	1.03
3:H:27:ALA:HB1	3:H:28:SER:HA	1.39	1.03
1:G:117:GLN:HE21	1:G:203:GLN:HG3	1.21	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:363:PRO:O	1:G:469:ARG:NH1	1.99	0.95
1:G:117:GLN:NE2	1:G:203:GLN:HG3	1.83	0.93

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	G	324/353 (92%)	246 (76%)	46 (14%)	32 (10%)	1	3
2	N	24/28 (86%)	21 (88%)	2 (8%)	1 (4%)	3	18
3	H	215/229 (94%)	168 (78%)	24 (11%)	23 (11%)	0	2
4	L	210/221 (95%)	171 (81%)	28 (13%)	11 (5%)	2	13
All	All	773/831 (93%)	606 (78%)	100 (13%)	67 (9%)	1	4

5 of 67 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	60	ALA
1	G	94	ASN
1	G	114	GLN
1	G	204	ALA
1	G	205	CYS

5.3.2 Protein sidechains [i](#)

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	G	297/310 (96%)	271 (91%)	26 (9%)	12	41
2	N	20/20 (100%)	19 (95%)	1 (5%)	30	68
3	H	188/197 (95%)	166 (88%)	22 (12%)	7	25
4	L	180/188 (96%)	159 (88%)	21 (12%)	7	25
All	All	685/715 (96%)	615 (90%)	70 (10%)	9	32

5 of 70 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	H	35(B)	SER
3	H	96	PRO
4	L	165	SER
3	H	38	ARG
3	H	71	VAL

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

Mol	Chain	Res	Type
1	G	117	GLN
1	G	258	GLN
1	G	462	ASN
3	H	197	ASN
4	L	1	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	DPR	N	21	2	5,7,8	0.58	0	7,8,10	1.38	1 (14%)
2	U2X	N	23	2	18,20,21	1.14	1 (5%)	23,25,27	1.21	1 (4%)

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	DPR	N	21	2	-	0/0/9/11	0/1/1/1
2	U2X	N	23	2	-	0/9/19/21	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	N	23	U2X	OH-CZ	3.38	1.45	1.37

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	23	U2X	O-C-CA	-4.48	113.71	125.72
2	N	21	DPR	CG-CB-CA	2.01	109.00	104.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

11 ligands 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
5	NAG	G	501	1	14,14,15	0.57	0	15,19,21	1.06	2 (13%)
5	NAG	G	502	1	14,14,15	0.38	0	15,19,21	1.63	1 (6%)
5	NAG	G	503	1	14,14,15	0.63	0	15,19,21	1.16	1 (6%)
5	NAG	G	504	1	14,14,15	0.43	0	15,19,21	1.46	3 (20%)
5	NAG	G	505	1	14,14,15	0.41	0	15,19,21	1.83	3 (20%)
5	NAG	G	506	1	14,14,15	0.42	0	15,19,21	0.74	0
5	NAG	G	507	1	14,14,15	0.38	0	15,19,21	1.01	1 (6%)
5	NAG	G	508	1	14,14,15	0.61	0	15,19,21	1.49	2 (13%)
5	NAG	G	509	1	14,14,15	1.67	2 (14%)	15,19,21	2.60	1 (6%)
5	NAG	G	510	1	14,14,15	0.61	0	15,19,21	1.49	2 (13%)
6	MPD	L	301	-	6,7,7	0.36	0	6,10,10	0.38	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
5	NAG	G	501	1	-	0/6/23/26	0/1/1/1
5	NAG	G	502	1	-	0/6/23/26	0/1/1/1
5	NAG	G	503	1	-	0/6/23/26	0/1/1/1
5	NAG	G	504	1	-	0/6/23/26	0/1/1/1
5	NAG	G	505	1	-	0/6/23/26	0/1/1/1
5	NAG	G	506	1	-	0/6/23/26	0/1/1/1
5	NAG	G	507	1	-	0/6/23/26	0/1/1/1
5	NAG	G	508	1	-	0/6/23/26	0/1/1/1
5	NAG	G	509	1	-	0/6/23/26	0/1/1/1
5	NAG	G	510	1	-	0/6/23/26	0/1/1/1
6	MPD	L	301	-	-	0/5/5/5	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	G	509	NAG	C2-N2	2.29	1.50	1.46
5	G	509	NAG	C1-C2	5.40	1.60	1.52

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	508	NAG	C1-O5-C5	-3.20	107.44	112.14
5	G	505	NAG	C4-C3-C2	-2.78	107.02	111.34
5	G	507	NAG	O5-C5-C4	-2.73	105.61	110.13
5	G	508	NAG	O5-C5-C4	-2.53	105.95	110.13
5	G	504	NAG	C2-N2-C7	2.03	125.75	123.11

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	G	332/353 (94%)	0.12	14 (4%) 40 22	31, 56, 116, 143	0
2	N	24/28 (85%)	-0.27	0 100 100	33, 43, 52, 62	0
3	H	218/229 (95%)	-0.19	0 100 100	29, 41, 59, 66	0
4	L	212/221 (95%)	-0.24	2 (0%) 85 69	33, 45, 60, 75	0
All	All	786/831 (94%)	-0.08	16 (2%) 68 46	29, 46, 91, 143	0

The worst 5 of 16 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	429	GLY	6.6
1	G	433	ALA	5.1
1	G	428	GLN	4.5
1	G	325	ASP	3.8
1	G	430	THR	3.5

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

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	U2X	N	23	19/20	0.95	0.20	-	32,35,39,39	0
2	DPR	N	21	7/8	0.94	0.19	-	42,43,47,51	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

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
5	NAG	G	510	14/15	0.66	0.34	5.11	79,87,94,96	0
6	MPD	L	301	8/8	0.88	0.23	2.72	54,59,67,70	0
5	NAG	G	501	14/15	0.89	0.20	2.23	51,55,56,58	0
5	NAG	G	509	14/15	0.79	0.30	1.63	76,79,86,86	0
5	NAG	G	505	14/15	0.90	0.21	0.20	47,49,61,63	0
5	NAG	G	506	14/15	0.92	0.23	0.11	46,51,54,55	0
5	NAG	G	503	14/15	0.94	0.17	-0.48	34,38,40,41	0
5	NAG	G	507	14/15	0.85	0.28	-	64,71,77,78	0
5	NAG	G	508	14/15	0.78	0.28	-	59,65,69,69	0
5	NAG	G	504	14/15	0.82	0.26	-	57,63,72,74	0
5	NAG	G	502	14/15	0.82	0.31	-	79,88,99,101	0

6.5 Other polymers [i](#)

There are no such residues in this entry.