



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:43 AM GMT

PDB ID : 2NY3
Title : HIV-1 gp120 Envelope Glycoprotein (K231C, T257S, E267C, S334A, S375W)
Complexed with CD4 and Antibody 17b
Authors : Zhou, T.; Xu, L.; Dey, B.; Hessel, A.J.; Van Ryk, D.; Xiang, S.H.; Yang, X.;
Zhang, M.Y.; Zwick, M.B.; Arthos, J.; Burton, D.R.; Dimitrov, D.S.; Sodroski,
J.; Wyatt, R.; Nabel, G.J.; Kwong, P.D.
Deposited on : 2006-11-20
Resolution : 2.00 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

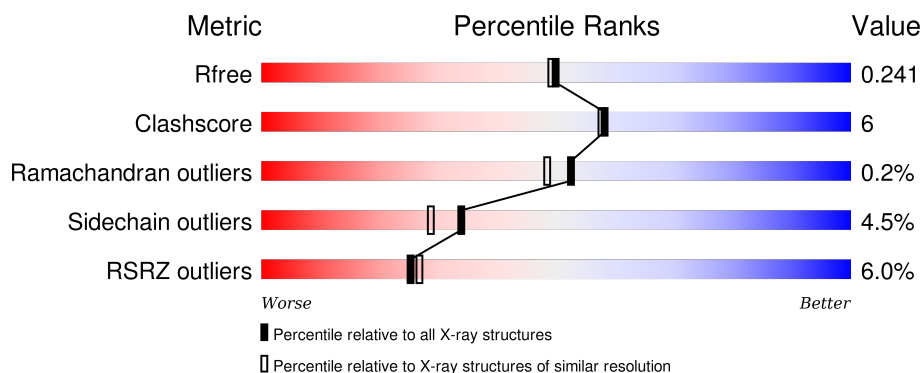
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	317	<div> <div>10%</div> <div>86%</div> <div>9%</div> <div>• •</div> </div>
2	B	184	<div> <div>6%</div> <div>78%</div> <div>18%</div> <div>• •</div> </div>
3	C	214	<div> <div>2%</div> <div>88%</div> <div>10%</div> <div>•</div> </div>
4	D	229	<div> <div>3%</div> <div>86%</div> <div>10%</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
5	NAG	A	730	-	-	-	X
5	NAG	A	734	-	-	-	X
6	SUC	D	4000	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7974 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 ENVELOPE GLYCOPROTEIN GP120.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	0	0
			2357	1480	411	444	22			

- Molecule 2 is a protein called T-cell surface glycoprotein CD4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	181	Total	C	N	O	S	0	0	0
			1412	885	247	276	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1000	MET	-	INSERTION	UNP P01730

- Molecule 3 is a protein called ANTIBODY 17B, LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	214	Total	C	N	O	S	0	0	0
			1646	1028	282	331	5			

- Molecule 4 is a protein called ANTIBODY 17B, HEAVY CHAIN.

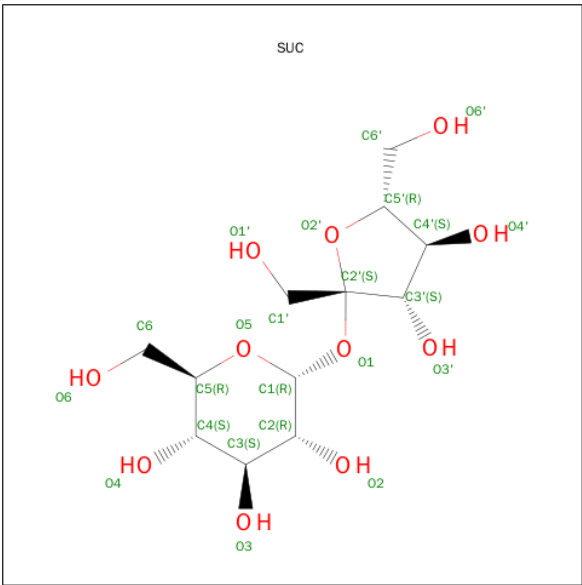
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	223	Total	C	N	O	S	0	2	0
			1689	1068	285	331	5			

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



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

- Molecule 6 is SUGAR (SUCROSE) (three-letter code: SUC) (formula: C₁₂H₂₂O₁₁).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	C	O	0	0
			23	12	11		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	215	Total	O	0	0
			215	215		
7	B	99	Total	O	0	0
			99	99		
7	C	162	Total	O	0	0
			162	162		
7	D	189	Total	O	0	0
			189	189		



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	71.74Å 88.20Å 197.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.11 – 2.00 44.10 – 2.00	Depositor EDS
% Data completeness (in resolution range)	91.6 (44.11-2.00) 91.6 (44.10-2.00)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.203 , 0.243 0.202 , 0.241	Depositor DCC
R_{free} test set	3926 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.1	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 78267 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7974	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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: SUC, NAG

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	0.33	1/2405 (0.0%)	0.48	1/3263 (0.0%)
2	B	0.29	0/1432	0.48	0/1930
3	C	0.35	1/1683 (0.1%)	0.51	0/2288
4	D	0.32	0/1740	0.50	0/2370
All	All	0.32	2/7260 (0.0%)	0.49	1/9851 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	2192	LYS	CE-NZ	6.20	1.64	1.49
1	A	491	ILE	C-O	5.04	1.32	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	452	LEU	CA-CB-CG	5.49	127.93	115.30

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	2357	0	2296	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1412	0	1441	31	0
3	C	1646	0	1590	12	0
4	D	1689	0	1650	23	0
5	A	182	0	169	3	0
6	D	23	0	22	1	0
7	A	215	0	0	7	0
7	B	99	0	0	12	0
7	C	162	0	0	1	0
7	D	189	0	0	6	0
All	All	7974	0	7168	86	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 6.

All (86) 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)
4:D:3215:HIS:HB3	4:D:3220:THR:HG22	1.33	1.05
2:B:1068:PRO:HA	7:B:316:HOH:O	1.57	1.04
1:A:350:ARG:HH11	1:A:350:ARG:HG3	1.33	0.93
2:B:1025:GLN:HG2	7:B:153:HOH:O	1.69	0.91
1:A:419:ARG:NH2	4:D:3103:GLU:OE2	2.06	0.89
4:D:3003:GLN:OE1	7:D:636:HOH:O	1.94	0.85
4:D:3065:GLN:NE2	7:D:36:HOH:O	2.11	0.84
1:A:350:ARG:HH11	1:A:350:ARG:CG	1.95	0.79
4:D:3046:GLU:OE2	4:D:3063[A]:HIS:NE2	2.15	0.79
2:B:1076:ILE:HD12	7:B:173:HOH:O	1.84	0.76
4:D:3112:ASN:ND2	7:D:506:HOH:O	2.20	0.75
4:D:3215:HIS:HB3	4:D:3220:THR:CG2	2.15	0.75
1:A:480:ARG:NH2	7:A:1093:HOH:O	2.21	0.73
2:B:1097:VAL:O	7:B:173:HOH:O	2.06	0.72
2:B:1027:HIS:CD2	7:B:501:HOH:O	2.41	0.71
4:D:3101:GLU:HA	7:D:506:HOH:O	1.90	0.70
2:B:1028:TRP:HZ2	7:B:316:HOH:O	1.76	0.69
5:A:734:NAG:O7	7:A:1163:HOH:O	2.08	0.69
4:D:3134:PRO:HD2	4:D:3220:THR:HG21	1.76	0.67
1:A:123:THR:HG21	1:A:429:LYS:HE3	1.76	0.67
4:D:3110:ASP:HB2	7:D:506:HOH:O	1.95	0.66
2:B:1029:LYS:HG2	7:B:501:HOH:O	1.95	0.65
1:A:119:CYS:N	1:A:205:CYS:SG	2.70	0.65
2:B:1164:ASN:O	2:B:1166:LYS:HD2	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1029:LYS:CG	7:B:501:HOH:O	2.46	0.62
2:B:1120:SER:HA	7:B:173:HOH:O	2.01	0.61
4:D:3063[A]:HIS:CD2	4:D:3064:LEU:HD13	2.37	0.59
4:D:3063[A]:HIS:HD2	4:D:3064:LEU:HD13	1.68	0.59
1:A:294:ILE:HD12	1:A:333:ILE:HD11	1.85	0.58
4:D:3063[A]:HIS:CE1	6:D:4000:SUC:O6'	2.58	0.56
3:C:2188:TYR:CZ	3:C:2213:ARG:HG3	2.40	0.56
2:B:1083:ILE:HG13	2:B:1092:GLU:HG3	1.87	0.56
2:B:1157:TRP:HB3	7:B:527:HOH:O	2.05	0.56
5:A:762:NAG:O4	7:A:1141:HOH:O	2.07	0.56
1:A:295:ASN:HD22	5:A:795:NAG:H83	1.72	0.55
1:A:122:LEU:HD22	1:A:200:VAL:HG22	1.88	0.55
3:C:2171:LYS:HA	3:C:2171:LYS:HE3	1.89	0.54
2:B:1142:LYS:HA	2:B:1142:LYS:NZ	2.24	0.53
3:C:2029:VAL:HG11	3:C:2090:GLN:HG2	1.91	0.52
4:D:3066:GLY:HA3	7:D:254:HOH:O	2.09	0.52
4:D:3203:SER:HG	4:D:3209:TYR:HH	1.56	0.52
3:C:2094:TRP:HA	3:C:2095:PRO:C	2.29	0.52
3:C:2185:LYS:HE2	3:C:2189:GLU:OE2	2.10	0.52
2:B:1084:CYS:HB3	2:B:1091:GLU:HG3	1.93	0.51
2:B:1142:LYS:HA	2:B:1142:LYS:HZ2	1.75	0.51
2:B:1120:SER:OG	2:B:1142:LYS:HE2	2.10	0.51
2:B:1138:ILE:HG21	2:B:1146:VAL:HG22	1.93	0.50
3:C:2029:VAL:HG11	3:C:2090:GLN:CG	2.41	0.50
4:D:3038:ARG:HD2	4:D:3046:GLU:OE1	2.12	0.50
3:C:2037:GLN:HB2	3:C:2047:LEU:HD11	1.93	0.49
3:C:2054:ARG:HG2	3:C:2058:VAL:HB	1.94	0.49
3:C:2004:MET:SD	3:C:2090:GLN:HB2	2.53	0.49
1:A:394:THR:HG22	1:A:397:ASN:HB2	1.95	0.48
2:B:1029:LYS:HG3	7:B:501:HOH:O	2.13	0.47
1:A:385:CYS:SG	1:A:417:PRO:O	2.73	0.47
4:D:3081:LEU:C	4:D:3081:LEU:HD13	2.35	0.47
3:C:2122:PRO:HD3	3:C:2134:VAL:HG22	1.98	0.46
3:C:2105:ARG:NH1	7:C:168:HOH:O	2.49	0.45
2:B:1012:VAL:HB	2:B:1074:LEU:HD11	1.97	0.45
2:B:1094:GLN:HE21	2:B:1096:LEU:HD21	1.82	0.45
1:A:273:ARG:NH2	1:A:287:GLN:OE1	2.49	0.45
2:B:1021:LYS:HE3	2:B:1064:GLN:OE1	2.17	0.44
1:A:350:ARG:NH1	1:A:350:ARG:CG	2.62	0.44
2:B:1181:LYS:HB3	2:B:1181:LYS:HE3	1.84	0.44
4:D:3038:ARG:HG3	4:D:3046:GLU:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:3105:ASP:OD1	4:D:3105:ASP:N	2.51	0.44
1:A:371:ILE:HG12	2:B:1043:PHE:HB3	2.00	0.43
1:A:480:ARG:NE	7:A:1155:HOH:O	2.44	0.43
4:D:3165:VAL:CG1	4:D:3193:LEU:HD21	2.49	0.42
2:B:1162:LEU:HD23	2:B:1163:GLN:N	2.34	0.42
4:D:3216:LYS:N	4:D:3217:PRO:CD	2.83	0.42
1:A:269:GLU:HB2	7:A:1161:HOH:O	2.20	0.42
1:A:442:GLN:HG2	1:A:444:ARG:HH12	1.85	0.42
2:B:1027:HIS:HD2	7:B:501:HOH:O	1.93	0.41
2:B:1138:ILE:HD12	2:B:1144:LEU:HB3	2.01	0.41
1:A:444:ARG:HG3	7:A:1074:HOH:O	2.20	0.41
1:A:96:TRP:CD1	1:A:236:THR:HG21	2.56	0.41
2:B:1118:LEU:O	2:B:1142:LYS:HE3	2.21	0.41
1:A:85:VAL:HA	1:A:243:SER:HB3	2.03	0.41
2:B:1164:ASN:O	2:B:1165:GLN:HB2	2.20	0.40
2:B:1077:GLU:CD	2:B:1077:GLU:H	2.24	0.40
2:B:1118:LEU:HD23	2:B:1142:LYS:NZ	2.37	0.40
3:C:2013:VAL:CG1	3:C:2078:LEU:HD12	2.51	0.40
4:D:3193:LEU:HD12	4:D:3193:LEU:C	2.42	0.40
2:B:1021:LYS:NZ	2:B:1064:GLN:HG3	2.36	0.40
7:A:1106:HOH:O	4:D:3108:GLU:HB2	2.21	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	301/317 (95%)	283 (94%)	17 (6%)	1 (0%)	46	41
2	B	179/184 (97%)	174 (97%)	5 (3%)	0	100	100
3	C	212/214 (99%)	205 (97%)	7 (3%)	0	100	100
4	D	221/229 (96%)	216 (98%)	4 (2%)	1 (0%)	34	26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	913/944 (97%)	878 (96%)	33 (4%)	2 (0%)	52	48

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	194	GLY
4	D	3228	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	269/280 (96%)	260 (97%)	9 (3%)	45	43
2	B	164/166 (99%)	153 (93%)	11 (7%)	20	14
3	C	184/184 (100%)	176 (96%)	8 (4%)	35	30
4	D	189/193 (98%)	181 (96%)	8 (4%)	36	31
All	All	806/823 (98%)	770 (96%)	36 (4%)	34	29

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	A	205	CYS
1	A	267	CYS
1	A	339	ASN
1	A	350	ARG
1	A	357	LYS
1	A	371	ILE
1	A	385	CYS
1	A	452	LEU
2	B	1002	LYS
2	B	1025	GLN
2	B	1027	HIS
2	B	1037	LEU

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Mol	Chain	Res	Type
2	B	1064	GLN
2	B	1069	LEU
2	B	1091	GLU
2	B	1100	LEU
2	B	1109	LEU
2	B	1142	LYS
2	B	1149	LEU
3	C	2054	ARG
3	C	2090	GLN
3	C	2107	GLU
3	C	2127	LEU
3	C	2131	THR
3	C	2156	LEU
3	C	2171	LYS
3	C	2183	LEU
4	D	3038	ARG
4	D	3055	LEU
4	D	3064	LEU
4	D	3083	LEU
4	D	3089	ASP
4	D	3165	VAL
4	D	3203	SER
4	D	3220	THR

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

Mol	Chain	Res	Type
1	A	98	ASN
1	A	355	ASN
2	B	1094	GLN
3	C	2102	GLN
4	D	3003	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

14 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	A	588	1	14,14,15	0.50	0	15,19,21	0.69	0
5	NAG	A	697	1	14,14,15	0.43	0	15,19,21	0.88	1 (6%)
5	NAG	A	730	1	14,14,15	0.59	0	15,19,21	0.92	1 (6%)
5	NAG	A	734	1	14,14,15	0.50	0	15,19,21	2.02	3 (20%)
5	NAG	A	741	1	14,14,15	0.53	0	15,19,21	0.89	1 (6%)
5	NAG	A	762	1	14,14,15	0.54	0	15,19,21	0.82	0
5	NAG	A	776	1	14,14,15	0.50	0	15,19,21	0.83	0
5	NAG	A	789	1	14,14,15	0.53	0	15,19,21	0.73	0
5	NAG	A	795	1	14,14,15	0.50	0	15,19,21	0.85	0
5	NAG	A	839	1	14,14,15	0.69	0	15,19,21	1.24	1 (6%)
5	NAG	A	886	1	14,14,15	0.48	0	15,19,21	0.61	0
5	NAG	A	892	1	14,14,15	0.55	0	15,19,21	0.71	0
5	NAG	A	948	1	14,14,15	0.47	0	15,19,21	0.83	0
6	SUC	D	4000	-	24,24,24	0.41	0	36,36,36	0.77	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	A	588	1	-	0/6/23/26	0/1/1/1
5	NAG	A	697	1	-	0/6/23/26	0/1/1/1
5	NAG	A	730	1	-	0/6/23/26	0/1/1/1
5	NAG	A	734	1	-	0/6/23/26	0/1/1/1
5	NAG	A	741	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	762	1	-	0/6/23/26	0/1/1/1
5	NAG	A	776	1	-	0/6/23/26	0/1/1/1
5	NAG	A	789	1	-	0/6/23/26	0/1/1/1
5	NAG	A	795	1	-	0/6/23/26	0/1/1/1
5	NAG	A	839	1	-	0/6/23/26	0/1/1/1
5	NAG	A	886	1	-	0/6/23/26	0/1/1/1
5	NAG	A	892	1	-	0/6/23/26	0/1/1/1
5	NAG	A	948	1	-	0/6/23/26	0/1/1/1
6	SUC	D	4000	-	-	0/12/51/51	0/2/2/2

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	734	NAG	C4-C3-C2	-3.24	106.19	111.23
5	A	734	NAG	C3-C4-C5	-2.41	106.00	110.20
5	A	697	NAG	C2-N2-C7	-2.16	120.26	123.04
5	A	741	NAG	C4-C3-C2	2.00	114.34	111.23
5	A	730	NAG	C4-C3-C2	2.33	114.86	111.23
5	A	839	NAG	C4-C3-C2	3.83	117.18	111.23
5	A	734	NAG	C1-O5-C5	6.34	120.30	112.25

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
5	A	734	NAG	1	0
5	A	762	NAG	1	0
5	A	795	NAG	1	0
6	D	4000	SUC	1	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	305/317 (96%)	0.54	31 (10%) 9 9	14, 30, 65, 95	0
2	B	181/184 (98%)	0.33	11 (6%) 25 26	16, 35, 59, 70	0
3	C	214/214 (100%)	-0.04	5 (2%) 64 64	12, 27, 53, 64	0
4	D	223/229 (97%)	0.15	8 (3%) 46 48	13, 23, 55, 77	0
All	All	923/944 (97%)	0.27	55 (5%) 25 27	12, 29, 59, 95	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	410	GLY	7.4
1	A	194	GLY	6.1
1	A	128	GLY	5.4
1	A	129	ALA	5.2
1	A	462	ASN	5.1
1	A	398	SER	5.1
1	A	385	CYS	4.5
2	B	1105	ASP	4.5
4	D	3148	GLY	4.2
1	A	88	ASN	3.8
2	B	1001	LYS	3.7
1	A	126	CYS	3.6
1	A	84	VAL	3.5
1	A	87	VAL	3.4
1	A	195	SER	3.4
1	A	205	CYS	3.3
4	D	3229	LYS	3.2
1	A	196	CYS	3.2
1	A	228	CYS	3.2
1	A	463	ASN	3.1
1	A	89	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	461	SER	3.0
1	A	83	GLU	3.0
1	A	127	VAL	3.0
3	C	2196	CYS	2.9
4	D	3228	PRO	2.9
1	A	85	VAL	2.8
1	A	86	LEU	2.7
2	B	1141	GLY	2.7
3	C	2001	ASP	2.7
2	B	1181	LYS	2.7
1	A	231	CYS	2.7
1	A	356	ASN	2.6
2	B	1106	THR	2.6
3	C	2214	GLY	2.6
2	B	1125	SER	2.6
1	A	464	GLU	2.4
2	B	1088	ASP	2.4
4	D	3001	GLU	2.4
1	A	210	PHE	2.4
2	B	1180	GLN	2.4
4	D	3149	GLY	2.4
2	B	1091	GLU	2.3
1	A	396	PHE	2.3
4	D	3205	GLY	2.3
4	D	3227	GLU	2.3
2	B	1019	SER	2.2
3	C	2124	ASP	2.2
1	A	99	ASP	2.2
1	A	267	CYS	2.1
4	D	3220	THR	2.1
1	A	397	ASN	2.1
3	C	2007	SER	2.0
1	A	119	CYS	2.0
2	B	1053	ASP	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 [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	A	730	14/15	0.65	0.43	7.14	54,56,56,57	0
5	NAG	A	734	14/15	0.71	0.33	5.98	49,50,51,51	0
6	SUC	D	4000	23/23	0.73	0.35	5.74	54,54,55,56	0
5	NAG	A	789	14/15	0.93	0.19	0.91	38,39,40,40	0
5	NAG	A	948	14/15	0.81	0.21	0.64	40,42,43,43	0
5	NAG	A	795	14/15	0.89	0.15	-0.03	31,33,34,34	0
5	NAG	A	762	14/15	0.91	0.15	-0.08	28,32,34,35	0
5	NAG	A	776	14/15	0.92	0.12	-0.56	36,38,39,39	0
5	NAG	A	697	14/15	0.55	0.52	-	53,55,55,56	0
5	NAG	A	588	14/15	0.55	0.43	-	60,61,62,62	0
5	NAG	A	741	14/15	0.62	0.44	-	52,53,54,54	0
5	NAG	A	839	14/15	0.67	0.43	-	48,50,51,51	0
5	NAG	A	886	14/15	0.90	0.12	-	37,38,39,39	0
5	NAG	A	892	14/15	0.79	0.34	-	42,44,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.