



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 11:55 pm GMT

PDB ID : 3MME  
Title : Structure and functional dissection of PG16, an antibody with broad and potent neutralization of HIV-1  
Authors : Pancera, M.; McLellan, J.; Zhou, T.; Zhu, J.; Kwong, P.  
Deposited on : 2010-04-19  
Resolution : 3.97 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](mailto: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.

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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

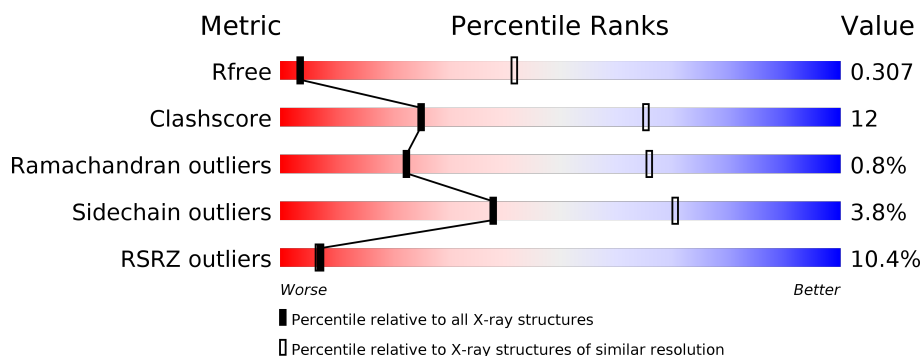
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.97 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	1066 (4.36-3.60)
Clashscore	112137	1163 (4.36-3.60)
Ramachandran outliers	110173	1119 (4.36-3.60)
Sidechain outliers	110143	1108 (4.36-3.60)
RSRZ outliers	101464	1078 (4.36-3.60)

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  $\geq 3$ , 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	238	
1	C	238	
1	H	238	
2	B	216	
2	D	216	
2	L	216	

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	D	570	-	-	X	-

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10040 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 PG16 HEAVY CHAIN FAB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	235	Total	C	N	O	S	0	0	0
			1802	1143	304	346	9			
1	A	235	Total	C	N	O	S	0	0	0
			1802	1143	304	346	9			
1	C	222	Total	C	N	O	S	0	0	0
			1684	1066	287	322	9			

- Molecule 2 is a protein called PG16 LIGHT CHAIN FAB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	211	Total	C	N	O	S	0	0	0
			1557	969	264	319	5			
2	B	211	Total	C	N	O	S	0	0	0
			1557	969	264	319	5			
2	D	211	Total	C	N	O	S	0	0	0
			1557	969	264	319	5			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	L	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



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

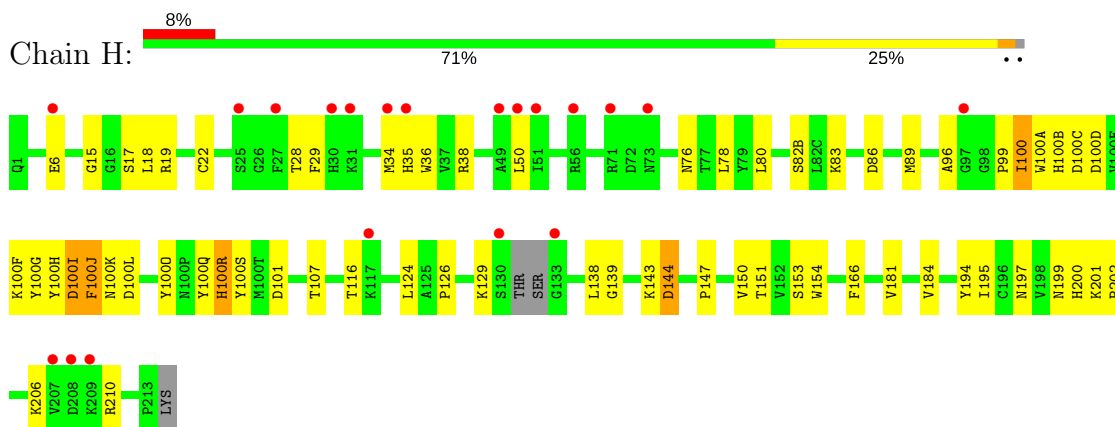
- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	2	Total	C	N	O	0	0
			28	16	2	10		

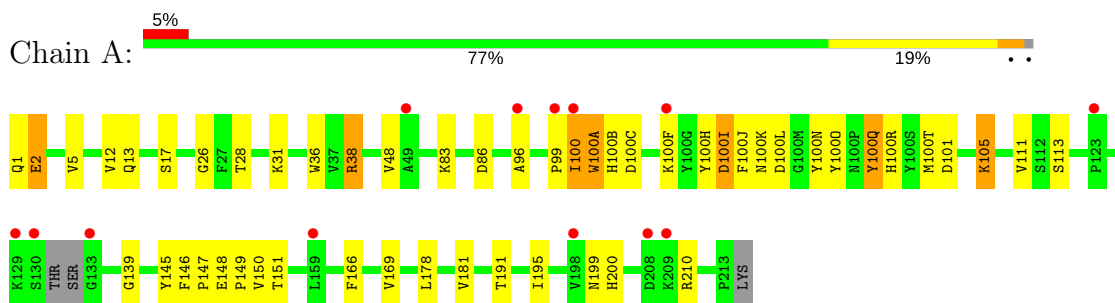
### 3 Residue-property plots [i](#)

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.

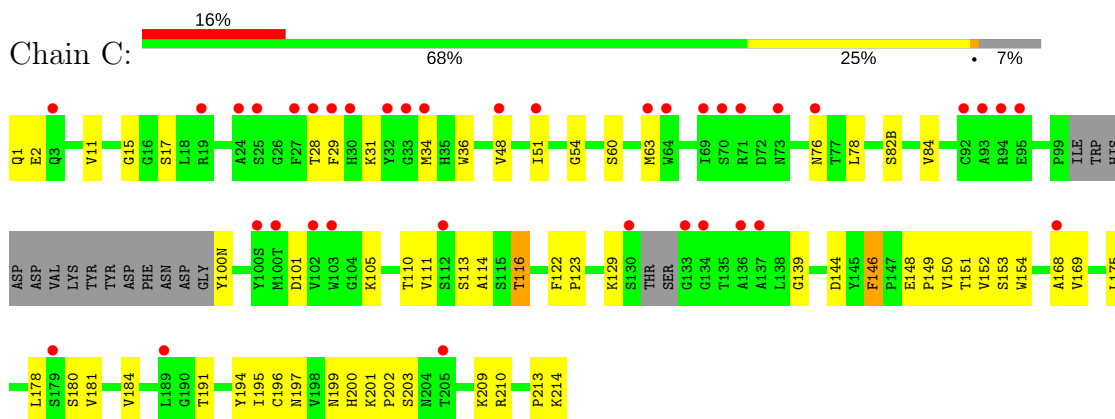
#### • Molecule 1: PG16 HEAVY CHAIN FAB



#### • Molecule 1: PG16 HEAVY CHAIN FAB

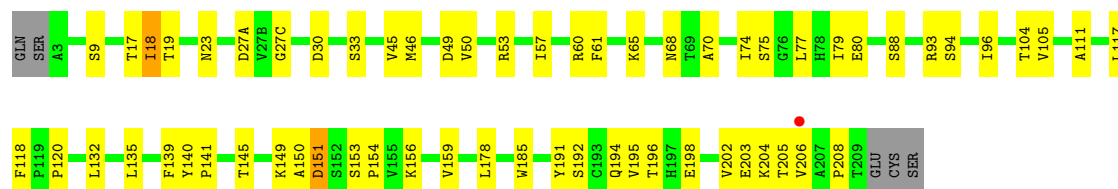


#### • Molecule 1: PG16 HEAVY CHAIN FAB



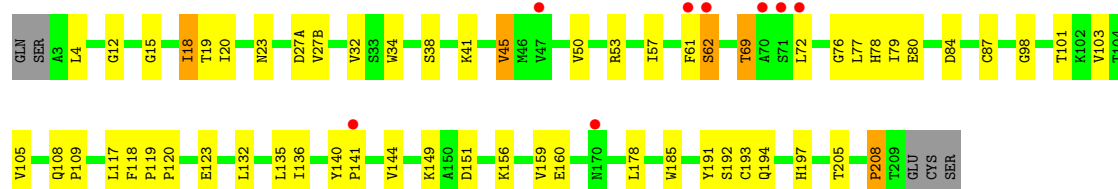
• Molecule 2: PG16 LIGHT CHAIN FAB

Chain L:  69% 28% ..




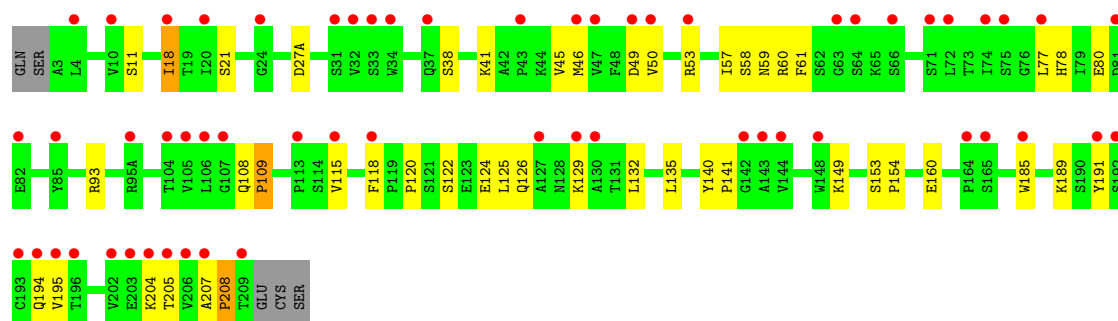
• Molecule 2: PG16 LIGHT CHAIN FAB

Chain B:  4% 70% 25% ..



• Molecule 2: PG16 LIGHT CHAIN FAB

Chain D:  27% 76% 20% ..



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	191.96Å 230.80Å 82.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.45 – 3.97 49.45 – 3.97	Depositor EDS
% Data completeness (in resolution range)	91.1 (49.45-3.97) 91.2 (49.45-3.97)	Depositor EDS
$R_{merge}$	0.24	Depositor
$R_{sym}$	0.20	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.73 (at 4.00Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, $R_{free}$	0.256 , 0.317 0.238 , 0.307	Depositor DCC
$R_{free}$ test set	738 reflections (5.01%)	DCC
Wilson B-factor (Å <sup>2</sup> )	112.9	Xtriage
Anisotropy	0.270	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 138.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.86	EDS
Total number of atoms	10040	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	176.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.75% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, 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	0.28	0/1852	0.44	0/2517
1	C	0.25	0/1727	0.43	0/2343
1	H	0.28	0/1852	0.44	0/2517
2	B	0.27	0/1594	0.44	0/2171
2	D	0.27	0/1594	0.43	0/2171
2	L	0.26	0/1594	0.45	0/2171
All	All	0.27	0/10213	0.44	0/13890

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 [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	1802	0	1730	46	0
1	C	1684	0	1640	41	0
1	H	1802	0	1730	48	0
2	B	1557	0	1507	38	0
2	D	1557	0	1507	40	0
2	L	1557	0	1507	42	0
3	L	39	0	34	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	14	0	13	1	0
5	D	28	0	25	8	0
All	All	10040	0	9693	238	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 12.

All (238) 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)
2:L:104:THR:HG21	2:L:141:PRO:HB3	1.46	0.97
1:H:181:VAL:HG11	2:L:135:LEU:HD13	1.53	0.90
2:D:38:SER:HB2	2:D:41:LYS:HD2	1.53	0.88
1:A:1:GLN:HG3	1:A:2:GLU:H	1.39	0.88
1:H:116:THR:HG22	1:H:147:PRO:HD3	1.58	0.83
1:A:1:GLN:HG3	1:A:2:GLU:HG2	1.61	0.81
1:H:195:ILE:HG12	1:H:210:ARG:HG2	1.63	0.80
1:C:181:VAL:HG11	2:D:135:LEU:HD13	1.63	0.80
2:D:21:SER:CB	5:D:570:NAG:HN2	1.96	0.78
1:H:150:VAL:HG22	1:H:200:HIS:HD2	1.50	0.77
2:L:18:ILE:HD12	2:L:77:LEU:HD11	1.69	0.75
1:A:100(I):ASP:HA	1:A:100(O):TYR:CZ	2.22	0.74
3:L:571:NAG:H3	3:L:572:BMA:H2	1.71	0.72
1:A:101:ASP:HA	2:B:45:VAL:HG21	1.72	0.69
1:H:101:ASP:HA	2:L:45:VAL:HG21	1.74	0.69
1:C:213:PRO:CB	1:C:214:LYS:HG3	2.22	0.69
2:D:21:SER:OG	5:D:570:NAG:N2	2.27	0.68
2:D:21:SER:CB	5:D:570:NAG:N2	2.57	0.67
1:A:100(A):TRP:HB3	1:A:100(F):LYS:HG2	1.77	0.66
1:A:169:VAL:HG21	2:B:160:GLU:HB3	1.78	0.66
2:L:156:LYS:NZ	2:D:189:LYS:HG3	2.10	0.66
2:D:21:SER:HB2	5:D:570:NAG:HN2	1.62	0.65
2:B:38:SER:HB2	2:B:41:LYS:HD2	1.77	0.65
2:D:21:SER:CB	5:D:570:NAG:C7	2.75	0.64
1:H:100(J):PHE:O	1:H:100(K):ASN:HB2	1.96	0.64
1:C:116:THR:HG22	1:C:203:SER:HB3	1.79	0.64
1:C:15:GLY:HA2	1:C:82(B):SER:HA	1.80	0.64
1:A:139:GLY:HA3	1:A:181:VAL:HG12	1.80	0.63
1:H:150:VAL:HG22	1:H:200:HIS:CD2	2.31	0.63
1:C:101:ASP:HA	2:D:45:VAL:HG21	1.81	0.62
1:A:195:ILE:HG12	1:A:210:ARG:HG2	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1:GLN:HG2	1:C:2:GLU:H	1.64	0.61
2:L:77:LEU:HD13	2:L:105:VAL:HG22	1.81	0.61
1:H:6:GLU:HG3	1:H:107:THR:HB	1.83	0.61
1:H:153:SER:OG	1:H:197:ASN:HB2	2.00	0.61
2:D:53:ARG:HD3	2:D:61:PHE:O	2.02	0.60
1:H:184:VAL:HG11	1:H:194:TYR:CE1	2.36	0.60
2:B:144:VAL:HG12	2:B:197:HIS:HB2	1.83	0.60
2:L:117:LEU:HG	2:L:206:VAL:HG13	1.83	0.59
1:C:36:TRP:O	1:C:48:VAL:HB	2.03	0.59
1:C:144:ASP:HA	1:C:175:LEU:HB3	1.84	0.59
1:C:181:VAL:HG11	2:D:135:LEU:CD1	2.31	0.59
2:B:53:ARG:HD3	2:B:61:PHE:O	2.03	0.58
2:D:21:SER:HB3	5:D:570:NAG:C7	2.33	0.58
2:L:159:VAL:HG22	2:L:178:LEU:HD13	1.85	0.58
1:H:151:THR:OG1	1:H:199:ASN:HB3	2.04	0.57
1:A:151:THR:OG1	1:A:199:ASN:HB3	2.04	0.57
1:A:96:ALA:O	1:A:100(Q):TYR:HB2	2.03	0.57
1:A:100(A):TRP:HD1	1:A:100(A):TRP:H	1.53	0.57
2:D:21:SER:OG	5:D:570:NAG:C7	2.53	0.56
1:H:36:TRP:CE2	1:H:80:LEU:HB2	2.40	0.56
1:H:100(K):ASN:O	1:H:100(L):ASP:HB2	2.05	0.56
1:C:34:MET:HB3	1:C:78:LEU:HD22	1.86	0.56
1:H:126:PRO:HG3	1:H:138:LEU:HD23	1.87	0.56
1:C:213:PRO:HB2	1:C:214:LYS:HG3	1.87	0.56
2:L:53:ARG:HD3	2:L:61:PHE:O	2.06	0.55
1:C:84:VAL:HA	1:C:111:VAL:HG13	1.87	0.55
1:C:123:PRO:HG3	1:C:209:LYS:HD2	1.89	0.55
1:A:150:VAL:HG22	1:A:200:HIS:HD2	1.72	0.55
2:L:145:THR:HB	2:L:196:THR:HB	1.88	0.55
1:H:38:ARG:HA	1:H:89:MET:O	2.06	0.54
1:C:184:VAL:HG11	1:C:194:TYR:CE1	2.42	0.54
1:H:6:GLU:HG2	1:H:107:THR:H	1.72	0.54
1:A:12:VAL:HG23	1:A:111:VAL:HG22	1.89	0.54
1:A:150:VAL:HG22	1:A:200:HIS:CD2	2.44	0.53
1:C:213:PRO:HB3	1:C:214:LYS:HG3	1.90	0.53
1:A:99:PRO:HG3	1:A:100(H):TYR:CE1	2.43	0.53
2:B:120:PRO:HD3	2:B:132:LEU:CD2	2.39	0.53
2:B:18:ILE:HD12	2:B:77:LEU:HD11	1.91	0.53
2:B:156:LYS:H	2:D:93:ARG:HH12	1.57	0.53
1:A:36:TRP:O	1:A:48:VAL:HB	2.09	0.53
2:B:4:LEU:HB2	2:B:98:GLY:HA2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:200:HIS:CE1	1:C:202:PRO:HB2	2.44	0.53
1:H:18:LEU:HD12	1:H:19:ARG:H	1.74	0.53
2:L:140:TYR:HA	2:L:141:PRO:C	2.29	0.53
1:A:166:PHE:CE1	2:B:135:LEU:HD22	2.44	0.52
2:B:149:LYS:HB2	2:B:192:SER:HB2	1.92	0.52
1:C:150:VAL:HG22	1:C:200:HIS:HD2	1.74	0.52
1:H:15:GLY:HA2	1:H:82(B):SER:HA	1.92	0.52
1:C:150:VAL:HG22	1:C:200:HIS:CD2	2.44	0.52
2:L:120:PRO:HD3	2:L:132:LEU:CD2	2.39	0.52
1:A:100(A):TRP:CB	1:A:100(F):LYS:HG2	2.39	0.52
1:C:195:ILE:HG12	1:C:210:ARG:HG2	1.91	0.52
2:L:17:THR:HG23	2:L:75:SER:HA	1.90	0.52
1:H:166:PHE:CE1	2:L:135:LEU:HD22	2.45	0.52
1:H:34:MET:HB3	1:H:78:LEU:HD22	1.90	0.52
1:H:100:ILE:HD13	1:H:100(A):TRP:N	2.25	0.51
1:H:100(I):ASP:HB3	1:H:100(O):TYR:HE1	1.75	0.51
2:B:18:ILE:HG23	2:B:19:THR:N	2.26	0.51
1:C:169:VAL:HG21	2:D:160:GLU:HB3	1.93	0.51
2:L:104:THR:HG21	2:L:141:PRO:CB	2.29	0.51
2:L:149:LYS:HB2	2:L:192:SER:HB2	1.93	0.51
2:D:118:PHE:HE2	2:D:135:LEU:HD12	1.74	0.51
1:C:148:GLU:HB3	1:C:149:PRO:HA	1.92	0.51
1:H:199:ASN:ND2	1:H:206:LYS:HE2	2.25	0.50
1:H:143:LYS:O	1:H:144:ASP:HB2	2.11	0.50
1:A:100(I):ASP:HA	1:A:100(O):TYR:CE1	2.45	0.50
2:L:46:MET:HA	2:L:57:ILE:HD13	1.93	0.50
1:A:38:ARG:HH12	1:A:86:ASP:HA	1.77	0.50
2:B:159:VAL:HG22	2:B:178:LEU:HD13	1.92	0.50
2:D:58:SER:OG	2:D:60:ARG:HG3	2.12	0.49
1:A:5:VAL:HG23	1:A:5:VAL:O	2.12	0.49
1:C:153:SER:OG	1:C:197:ASN:HB2	2.12	0.49
1:A:100(J):PHE:O	1:A:100(K):ASN:HB2	2.13	0.49
2:D:78:HIS:HB3	2:D:80:GLU:OE2	2.13	0.49
1:A:178:LEU:HD12	1:A:178:LEU:C	2.33	0.49
2:L:185:TRP:CZ2	2:L:208:PRO:HA	2.47	0.49
1:H:100(S):TYR:CD1	2:L:45:VAL:HG11	2.48	0.49
3:L:571:NAG:C3	3:L:572:BMA:H2	2.39	0.49
2:B:120:PRO:HD3	2:B:132:LEU:HD23	1.93	0.49
1:A:1:GLN:HG3	1:A:2:GLU:N	2.17	0.49
2:B:4:LEU:HB2	2:B:98:GLY:CA	2.43	0.49
1:H:22:CYS:HB3	1:H:78:LEU:HB3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:VAL:HG11	1:C:180:SER:CB	2.43	0.49
2:L:117:LEU:HG	2:L:206:VAL:CG1	2.43	0.49
1:C:29:PHE:CD2	1:C:76:ASN:HA	2.48	0.48
2:D:78:HIS:HB2	2:D:80:GLU:HG2	1.95	0.48
1:A:100(A):TRP:HB3	1:A:100(F):LYS:HA	1.94	0.48
1:C:168:ALA:HB2	1:C:178:LEU:HD23	1.95	0.48
1:H:29:PHE:CD2	1:H:76:ASN:HA	2.48	0.48
2:L:18:ILE:HG23	2:L:19:THR:N	2.28	0.48
3:L:571:NAG:H81	2:D:149:LYS:NZ	2.28	0.48
1:H:100(B):HIS:HB2	1:H:100(G):TYR:HE2	1.79	0.48
1:A:38:ARG:NH1	1:A:86:ASP:HA	2.28	0.48
1:A:148:GLU:HB3	1:A:149:PRO:HA	1.96	0.48
2:D:18:ILE:HD12	2:D:77:LEU:HD11	1.95	0.48
1:A:100(A):TRP:N	1:A:100(A):TRP:CD1	2.82	0.47
3:L:571:NAG:O7	2:D:149:LYS:NZ	2.41	0.47
2:L:65:LYS:HA	2:L:70:ALA:HA	1.96	0.47
2:D:125:LEU:HD23	2:D:129:LYS:O	2.14	0.47
1:H:126:PRO:HG3	1:H:138:LEU:CD2	2.45	0.47
1:C:1:GLN:HG2	1:C:2:GLU:HG3	1.97	0.47
1:H:96:ALA:O	1:H:100(Q):TYR:HB2	2.14	0.47
2:L:60:ARG:HB3	2:L:75:SER:O	2.15	0.47
1:A:100:ILE:HG22	1:A:100(O):TYR:CE2	2.49	0.47
1:C:129:LYS:NZ	2:D:204:LYS:HE2	2.30	0.47
2:D:46:MET:HA	2:D:57:ILE:HD13	1.97	0.47
2:B:34:TRP:CZ3	2:B:87:CYS:HB3	2.50	0.46
2:L:23:ASN:OD1	3:L:570:NAG:N2	2.48	0.46
1:A:31:LYS:HD3	1:A:100(N):TYR:CE2	2.50	0.46
2:D:122:SER:O	2:D:126:GLN:HG2	2.16	0.46
2:B:62:SER:O	2:B:72:LEU:HD12	2.15	0.46
2:B:185:TRP:CZ2	2:B:208:PRO:HA	2.51	0.46
1:H:100(Q):TYR:O	1:H:100(R):HIS:HB2	2.16	0.46
2:D:153:SER:HA	2:D:154:PRO:HD3	1.80	0.46
2:L:45:VAL:O	2:L:57:ILE:HD11	2.16	0.46
2:L:150:ALA:O	2:L:151:ASP:HB2	2.16	0.45
2:D:132:LEU:HD21	2:D:185:TRP:CZ3	2.51	0.45
1:C:51:ILE:HG12	1:C:54:GLY:HA2	1.98	0.45
1:H:35:HIS:CE1	1:H:50:LEU:HD13	2.52	0.45
1:H:96:ALA:C	1:H:100(Q):TYR:HB2	2.37	0.45
2:L:93:ARG:O	2:L:94:SER:HB2	2.17	0.45
1:C:201:LYS:N	1:C:202:PRO:CD	2.79	0.45
2:D:120:PRO:HD3	2:D:132:LEU:CD2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:120:PRO:HD3	2:L:132:LEU:HD23	1.99	0.45
2:L:153:SER:HA	2:L:154:PRO:HD3	1.85	0.45
2:B:78:HIS:HB2	2:B:80:GLU:HG2	1.99	0.45
2:D:140:TYR:HA	2:D:141:PRO:C	2.36	0.45
1:H:6:GLU:CG	1:H:107:THR:H	2.30	0.44
1:H:153:SER:HG	1:H:197:ASN:HB2	1.80	0.44
1:H:100(A):TRP:HB3	1:H:100(F):LYS:HG2	2.00	0.44
1:A:100(Q):TYR:HD2	1:A:100(Q):TYR:HA	1.68	0.44
1:C:122:PHE:CE1	2:D:124:GLU:HA	2.52	0.44
1:H:150:VAL:CG2	1:H:200:HIS:HD2	2.26	0.44
2:B:18:ILE:CG2	2:B:19:THR:N	2.80	0.44
2:L:191:TYR:O	2:L:205:THR:HG23	2.17	0.44
1:A:146:PHE:HA	1:A:147:PRO:HA	1.83	0.44
2:B:118:PHE:HA	2:B:119:PRO:HD3	1.78	0.44
1:C:105:LYS:HB3	1:C:105:LYS:HE2	1.85	0.44
1:C:113:SER:HA	1:C:114:ALA:HA	1.54	0.44
1:A:105:LYS:HE2	1:A:105:LYS:HB3	1.79	0.43
1:A:100(A):TRP:CG	1:A:100(F):LYS:HG2	2.54	0.43
1:A:100:ILE:HD13	1:A:100:ILE:C	2.38	0.43
2:L:104:THR:CG2	2:L:141:PRO:HB3	2.32	0.43
1:A:101:ASP:HA	2:B:45:VAL:CG2	2.42	0.43
2:B:84:ASP:HA	2:B:101:THR:O	2.18	0.43
5:D:570:NAG:H83	5:D:570:NAG:H3	1.99	0.43
1:H:129:LYS:NZ	2:L:204:LYS:HE2	2.33	0.43
2:B:23:ASN:HA	2:B:69:THR:HG23	2.01	0.43
1:H:129:LYS:HZ3	2:L:204:LYS:HE2	1.83	0.43
2:B:15:GLY:C	2:B:76:GLY:HA2	2.38	0.43
1:H:100(J):PHE:O	1:H:100(K):ASN:CB	2.66	0.43
2:B:191:TYR:O	2:B:205:THR:HG23	2.19	0.43
2:B:23:ASN:OD1	4:B:570:NAG:O5	2.35	0.42
1:C:146:PHE:C	1:C:146:PHE:CD2	2.93	0.42
1:H:99:PRO:HG3	1:H:100(H):TYR:CZ	2.54	0.42
1:A:145:TYR:CE1	1:A:150:VAL:HG23	2.54	0.42
1:C:60:SER:HB3	1:C:63:MET:HG2	2.01	0.42
2:D:11:SER:C	2:D:18:ILE:HD11	2.40	0.42
2:L:141:PRO:HG2	2:L:198:GLU:OE2	2.19	0.42
3:L:571:NAG:H2	3:L:571:NAG:H83	1.77	0.42
1:A:181:VAL:HG11	2:B:135:LEU:CD1	2.50	0.42
1:A:1:GLN:C	1:A:26:GLY:HA3	2.39	0.42
2:B:20:ILE:HD11	2:B:103:VAL:CG2	2.50	0.42
2:D:108:GLN:HA	2:D:109:PRO:HD3	1.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:GLN:CG	1:A:2:GLU:N	2.81	0.42
1:C:151:THR:OG1	1:C:199:ASN:HB3	2.19	0.42
2:D:149:LYS:HD3	2:D:194:GLN:NE2	2.34	0.42
2:L:27(C):GLY:HA3	2:L:68:ASN:OD1	2.19	0.42
1:C:11:VAL:HA	1:C:110:THR:O	2.19	0.42
1:C:122:PHE:HA	1:C:123:PRO:HD3	1.83	0.42
2:D:191:TYR:O	2:D:205:THR:HG23	2.19	0.42
1:A:13:GLN:OE1	1:A:113:SER:HA	2.19	0.42
1:C:101:ASP:HA	2:D:45:VAL:CG2	2.50	0.42
1:C:139:GLY:HA2	1:C:154:TRP:CH2	2.54	0.42
1:H:201:LYS:HB2	1:H:202:PRO:HD3	2.02	0.42
1:H:35:HIS:CE1	1:H:50:LEU:HB2	2.56	0.41
2:L:194:GLN:HG2	2:L:203:GLU:HG3	2.02	0.41
2:B:27(B):VAL:HG13	2:B:32:VAL:HG21	2.03	0.41
2:B:45:VAL:O	2:B:57:ILE:HD11	2.20	0.41
2:D:207:ALA:HA	2:D:208:PRO:HD3	1.90	0.41
2:L:202:VAL:HG12	2:L:203:GLU:N	2.36	0.41
2:L:74:ILE:HG21	2:L:77:LEU:HD23	2.02	0.41
1:A:100(H):TYR:HD2	1:A:100(L):ASP:O	2.04	0.41
2:B:12:GLY:O	2:B:105:VAL:HA	2.21	0.41
1:C:1:GLN:HG2	1:C:2:GLU:N	2.33	0.41
2:D:49:ASP:O	2:D:50:VAL:HB	2.20	0.41
1:H:100(H):TYR:N	1:H:100(H):TYR:CD1	2.88	0.41
2:B:108:GLN:HB3	2:B:109:PRO:HD2	2.03	0.41
2:B:149:LYS:HD3	2:B:194:GLN:NE2	2.35	0.41
2:L:88:SER:HA	2:L:96:ILE:O	2.20	0.41
1:A:83:LYS:O	1:A:86:ASP:HB2	2.21	0.41
1:H:100(Q):TYR:O	1:H:100(R):HIS:CB	2.68	0.41
1:H:139:GLY:HA2	1:H:154:TRP:CH2	2.55	0.41
3:L:570:NAG:O7	3:L:570:NAG:C3	2.69	0.41
2:B:78:HIS:HB3	2:B:80:GLU:OE2	2.20	0.41
1:C:31:LYS:HD3	1:C:100(N):TYR:CZ	2.56	0.41
1:H:83:LYS:O	1:H:86:ASP:HB2	2.20	0.41
2:B:140:TYR:HA	2:B:141:PRO:C	2.41	0.40
2:B:117:LEU:HD22	2:B:193:CYS:HB2	2.03	0.40
1:A:100(B):HIS:O	1:A:100(C):ASP:HB2	2.21	0.40
1:A:100(T):MET:O	2:B:45:VAL:HG22	2.21	0.40
3:L:571:NAG:C8	2:D:149:LYS:NZ	2.85	0.40
1:H:124:LEU:HB3	2:L:118:PHE:CD1	2.56	0.40
1:A:100:ILE:HD13	1:A:100(A):TRP:N	2.37	0.40
2:D:60:ARG:CZ	2:D:78:HIS:CD2	3.05	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:111:ALA:O	2:L:139:PHE:HA	2.21	0.40
2:L:49:ASP:O	2:L:50:VAL:HB	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	231/238 (97%)	214 (93%)	16 (7%)	1 (0%)	38	77
1	C	216/238 (91%)	202 (94%)	13 (6%)	1 (0%)	32	73
1	H	231/238 (97%)	209 (90%)	19 (8%)	3 (1%)	14	57
2	B	209/216 (97%)	196 (94%)	10 (5%)	3 (1%)	13	55
2	D	209/216 (97%)	196 (94%)	11 (5%)	2 (1%)	18	61
2	L	209/216 (97%)	197 (94%)	11 (5%)	1 (0%)	32	73
All	All	1305/1362 (96%)	1214 (93%)	80 (6%)	11 (1%)	22	66

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	100(R)	HIS
1	A	100(R)	HIS
1	H	100(C)	ASP
1	H	144	ASP
2	D	208	PRO
2	B	151	ASP
2	D	109	PRO
2	L	151	ASP
1	C	146	PHE
2	B	208	PRO
2	B	50	VAL



### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	199/202 (98%)	189 (95%)	10 (5%)	28	64
1	C	187/202 (93%)	182 (97%)	5 (3%)	50	77
1	H	199/202 (98%)	193 (97%)	6 (3%)	46	75
2	B	178/183 (97%)	170 (96%)	8 (4%)	32	67
2	D	178/183 (97%)	173 (97%)	5 (3%)	49	76
2	L	178/183 (97%)	170 (96%)	8 (4%)	32	67
All	All	1119/1155 (97%)	1077 (96%)	42 (4%)	38	70

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	17	SER
1	H	28	THR
1	H	100	ILE
1	H	100(D)	ASP
1	H	100(I)	ASP
1	H	100(J)	PHE
2	L	9	SER
2	L	18	ILE
2	L	27(A)	ASP
2	L	30	ASP
2	L	33	SER
2	L	79	ILE
2	L	80	GLU
2	L	195	VAL
1	A	2	GLU
1	A	17	SER
1	A	28	THR
1	A	38	ARG
1	A	100	ILE
1	A	100(A)	TRP
1	A	100(I)	ASP
1	A	100(Q)	TYR

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Mol	Chain	Res	Type
1	A	105	LYS
1	A	191	THR
2	B	18	ILE
2	B	27(A)	ASP
2	B	45	VAL
2	B	62	SER
2	B	69	THR
2	B	79	ILE
2	B	123	GLU
2	B	136	ILE
1	C	17	SER
1	C	28	THR
1	C	116	THR
1	C	191	THR
1	C	196	CYS
2	D	18	ILE
2	D	27(A)	ASP
2	D	59	ASN
2	D	115	VAL
2	D	195	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

5 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	$\# Z  > 2$
5	NAG	D	570	2,5	14,14,15	0.67	0	15,19,21	1.31	2 (13%)
5	NAG	D	571	5	14,14,15	0.52	0	15,19,21	0.73	0
3	NAG	L	570	3,2	14,14,15	0.96	1 (7%)	15,19,21	1.44	1 (6%)
3	NAG	L	571	3	14,14,15	0.42	0	15,19,21	1.48	3 (20%)
3	BMA	L	572	3	11,11,12	0.62	0	13,15,17	0.74	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	D	570	2,5	-	0/6/23/26	0/1/1/1
5	NAG	D	571	5	-	0/6/23/26	0/1/1/1
3	NAG	L	570	3,2	-	0/6/23/26	0/1/1/1
3	NAG	L	571	3	-	0/6/23/26	0/1/1/1
3	BMA	L	572	3	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	570	NAG	C1-C2	2.73	1.56	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	571	NAG	C2-N2-C7	-2.73	118.96	122.94
3	L	571	NAG	C4-C3-C2	-2.73	107.02	111.02
5	D	570	NAG	C4-C3-C2	2.86	115.21	111.02
3	L	571	NAG	C1-O5-C5	3.30	116.72	112.17
5	D	570	NAG	C1-O5-C5	3.37	116.81	112.17
3	L	570	NAG	C4-C3-C2	3.46	116.08	111.02

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	570	NAG	8	0
3	L	570	NAG	2	0
3	L	571	NAG	6	0
3	L	572	BMA	2	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	$\# Z  > 2$
4	NAG	B	570	2	14,14,15	0.45	0	15,19,21	1.42	2 (13%)

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
4	NAG	B	570	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	570	NAG	C4-C3-C2	-2.23	107.75	111.02
4	B	570	NAG	C1-O5-C5	3.93	117.59	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	570	NAG	1	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	235/238 (98%)	0.41	13 (5%) 26 20	98, 149, 242, 366	0
1	C	222/238 (93%)	0.96	38 (17%) 2 3	133, 209, 279, 350	0
1	H	235/238 (98%)	0.63	20 (8%) 11 10	98, 154, 238, 327	0
2	B	211/216 (97%)	0.42	8 (3%) 41 32	90, 138, 188, 222	0
2	D	211/216 (97%)	1.44	58 (27%) 1 1	149, 231, 314, 375	0
2	L	211/216 (97%)	0.19	1 (0%) 90 86	94, 140, 190, 235	0
All	All	1325/1362 (97%)	0.67	138 (10%) 7 7	90, 164, 271, 375	0

All (138) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	133	GLY	7.8
2	D	142	GLY	7.7
2	D	206	VAL	5.8
1	C	93	ALA	5.6
1	C	29	PHE	5.6
1	C	27	PHE	5.4
1	C	100(T)	MET	5.4
1	C	34	MET	4.9
2	D	72	LEU	4.8
2	D	113	PRO	4.7
2	D	203	GLU	4.7
2	D	74	ILE	4.7
2	D	104	THR	4.7
1	A	130	SER	4.4
2	D	20	ILE	4.3
1	H	130	SER	4.3
2	D	47	VAL	4.2
2	D	107	GLY	4.2
2	D	105	VAL	4.2

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Mol	Chain	Res	Type	RSRZ
1	C	102	VAL	4.2
2	D	143	ALA	4.1
1	C	136	ALA	4.0
1	H	6	GLU	4.0
2	D	207	ALA	4.0
2	D	18	ILE	3.9
2	D	195	VAL	3.9
2	D	95(A)	ARG	3.9
1	H	51	ILE	3.9
1	C	28	THR	3.8
2	D	71	SER	3.8
2	D	191	TYR	3.8
2	D	129	LYS	3.7
2	D	164	PRO	3.6
2	D	34	TRP	3.5
2	D	148	TRP	3.5
1	H	73	ASN	3.5
2	D	209	THR	3.4
1	A	133	GLY	3.4
1	A	159	LEU	3.4
2	D	144	VAL	3.4
2	D	85	TYR	3.3
1	C	130	SER	3.3
1	A	208	ASP	3.2
1	A	100	ILE	3.2
2	D	33	SER	3.2
2	D	185	TRP	3.2
2	D	46	MET	3.1
1	C	95	GLU	3.1
2	B	61	PHE	3.1
2	D	82	GLU	3.0
2	D	24	GLY	3.0
1	C	94	ARG	3.0
1	C	71	ARG	2.9
1	C	69	ILE	2.9
2	B	62	SER	2.9
2	D	75	SER	2.9
1	C	64	TRP	2.9
1	C	33	GLY	2.9
1	H	34	MET	2.9
1	A	129	LYS	2.9
1	H	35	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	73	ASN	2.9
2	B	47	VAL	2.8
2	D	64	SER	2.8
2	D	204	LYS	2.8
1	A	99	PRO	2.7
1	C	137	ALA	2.7
1	C	179	SER	2.7
1	C	19	ARG	2.7
1	H	209	LYS	2.7
1	A	198	VAL	2.7
1	C	32	TYR	2.7
1	C	30	HIS	2.6
2	D	63	GLY	2.6
1	H	133	GLY	2.6
2	L	206	VAL	2.6
2	D	193	CYS	2.6
2	D	50	VAL	2.6
2	D	77	LEU	2.6
1	C	134	GLY	2.6
2	D	106	LEU	2.5
1	C	3	GLN	2.5
1	C	205	THR	2.5
1	C	24	ALA	2.5
1	A	123	PRO	2.5
1	C	48	VAL	2.5
1	H	31	LYS	2.4
1	A	209	LYS	2.4
2	D	194	GLN	2.4
1	H	30	HIS	2.4
1	H	71	ARG	2.4
2	B	170	ASN	2.4
2	D	205	THR	2.4
1	C	70	SER	2.4
1	H	207	VAL	2.4
2	D	192	SER	2.4
1	C	76	ASN	2.3
2	D	49	ASP	2.3
1	C	168	ALA	2.3
1	C	112	SER	2.3
2	D	53	ARG	2.3
2	B	70	ALA	2.3
2	B	141	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	71	SER	2.2
2	D	127	ALA	2.2
2	D	115	VAL	2.2
1	H	49	ALA	2.2
1	H	50	LEU	2.2
2	D	202	VAL	2.2
2	D	165	SER	2.2
2	D	10	VAL	2.2
1	H	117	LYS	2.2
1	C	189	LEU	2.2
1	H	56	ARG	2.2
2	D	43	PRO	2.2
1	A	100(F)	LYS	2.2
2	D	118	PHE	2.2
2	B	72	LEU	2.1
1	H	25	SER	2.1
2	D	32	VAL	2.1
1	A	96	ALA	2.1
1	C	92	CYS	2.1
1	C	100(S)	TYR	2.1
2	D	66	SER	2.1
1	C	103	TRP	2.1
2	D	81	ASP	2.1
2	D	196	THR	2.1
1	C	51	ILE	2.1
1	H	208	ASP	2.1
1	C	63	MET	2.1
2	D	37	GLN	2.1
1	H	27	PHE	2.0
1	C	25	SER	2.0
1	H	97	GLY	2.0
2	D	130	ALA	2.0
2	D	31	SER	2.0
1	A	49	ALA	2.0
2	D	4	LEU	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](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	NAG	L	570	14/15	0.82	0.28	-	207,252,286,298	0
5	NAG	D	570	14/15	0.68	0.21	-	312,354,392,414	0
5	NAG	D	571	14/15	0.78	0.20	-	262,355,376,399	0
3	NAG	L	571	14/15	0.75	0.35	-	173,280,307,318	0
3	BMA	L	572	11/12	0.59	0.38	-	271,315,341,345	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	NAG	B	570	14/15	0.60	0.37	-	310,347,360,360	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.