



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 04:19 PM GMT

PDB ID : 1FMX
Title : STRUCTURE OF NATIVE PROTEINASE A IN THE SPACE GROUP P21
Authors : Gustchina, A.; Li, M.; Phylip, L.H.; Lees, W.E.; Kay, J.; Wlodawer, A.
Deposited on : 2000-08-18
Resolution : 2.61 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

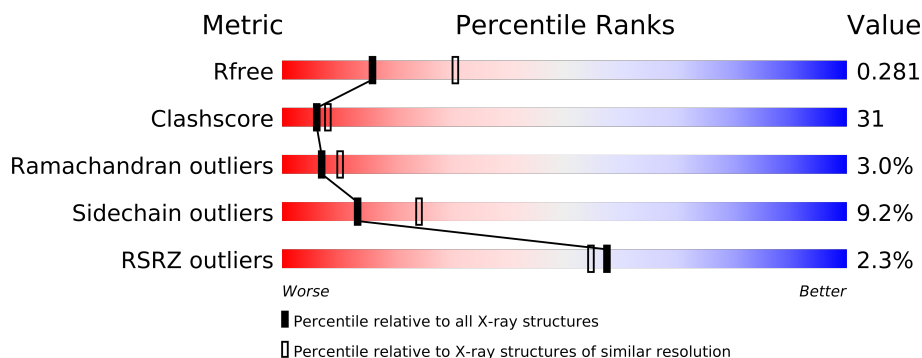
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.61 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}	66092	1940 (2.64-2.60)
Clashscore	79885	2404 (2.64-2.60)
Ramachandran outliers	78287	2360 (2.64-2.60)
Sidechain outliers	78261	2360 (2.64-2.60)
RSRZ outliers	66119	1939 (2.64-2.60)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	329	
1	B	329	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5288 atoms, of which 0 are hydrogen and 0 are deuterium.

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 SACCHAROPEPSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	51	0	0
			2498	1599	391	502	6			
1	B	325	Total	C	N	O	S	59	0	0
			2498	1599	391	502	6			

- Molecule 2 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	3	Total	C	N	O	4	0
			39	22	2	15		

- Molecule 3 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	2	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is water.

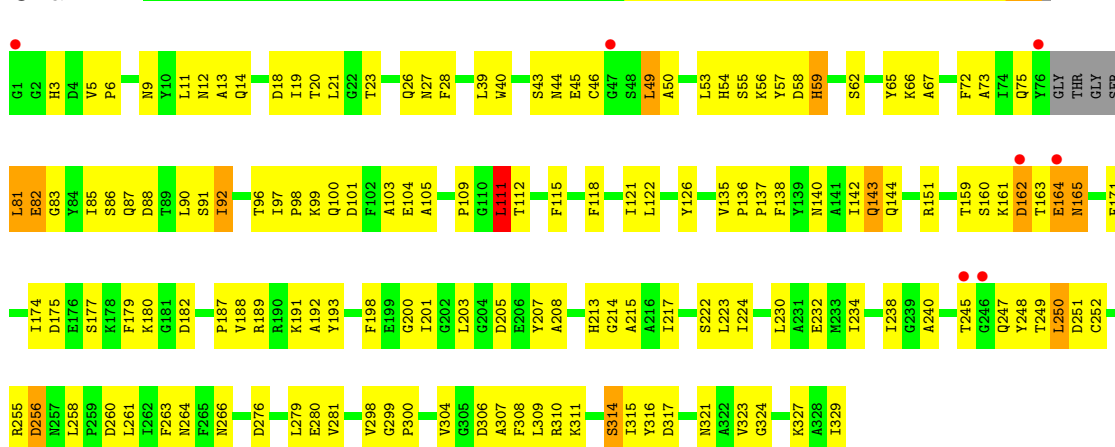
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	109	Total	O	0	0
			109	109		
5	B	88	Total	O	0	0
			88	88		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

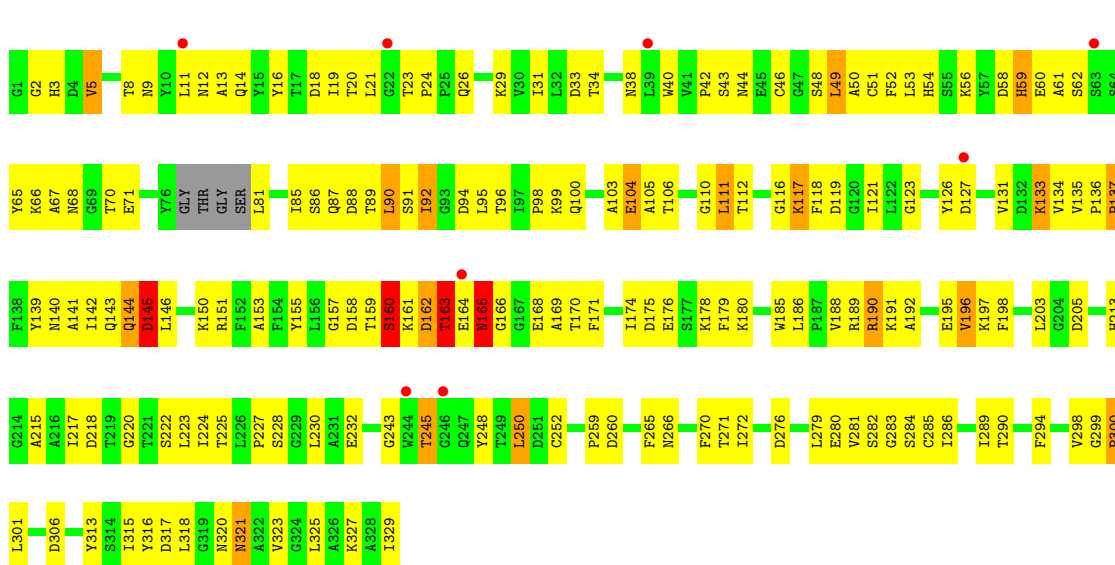
• Molecule 1: SACCHAROPEPSIN

Chain A:



• Molecule 1: SACCHAROPEPSIN

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.96Å 49.07Å 94.69Å 90.00° 96.50° 90.00°	Depositor
Resolution (Å)	24.99 – 2.61 24.99 – 2.61	Depositor EDS
% Data completeness (in resolution range)	91.3 (24.99-2.61) 94.4 (24.99-2.61)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.60Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.205 , 0.279 0.207 , 0.281	Depositor DCC
R_{free} test set	1080 reflections (5.15%)	DCC
Wilson B-factor (Å ²)	52.9	Xtriage
Anisotropy	0.164	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 53.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 22963 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5288	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.01% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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	0.67	0/2559	0.90	1/3479 (0.0%)
1	B	0.68	1/2559 (0.0%)	0.94	5/3479 (0.1%)
All	All	0.68	1/5118 (0.0%)	0.92	6/6958 (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	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	196	VAL	CB-CG1	5.58	1.64	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	162	ASP	CA-C-N	-10.47	94.17	117.20
1	B	165	ASN	N-CA-C	5.73	126.48	111.00
1	A	111	LEU	CA-CB-CG	5.61	128.20	115.30
1	B	162	ASP	C-N-CA	5.56	135.60	121.70
1	B	162	ASP	O-C-N	5.44	131.41	122.70
1	B	162	ASP	CA-C-O	5.04	130.68	120.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	193	TYR	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2498	0	2374	133	0
1	B	2498	0	2374	163	0
2	A	39	0	34	0	0
3	B	28	0	25	4	0
4	A	14	0	13	0	0
4	B	14	0	13	0	0
5	A	109	0	0	18	0
5	B	88	0	0	24	0
All	All	5288	0	4833	297	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 31.

All (297) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:121:ILE:HG22	5:B:382:HOH:O	1.64	0.96
1:B:134:VAL:HG12	5:B:408:HOH:O	1.64	0.95
1:A:234:ILE:HD13	1:A:261:LEU:HD11	1.49	0.95
1:B:5:VAL:HG21	1:B:92:ILE:HD12	1.50	0.94
1:B:44:ASN:ND2	1:B:58:ASP:HA	1.83	0.93
1:B:70:THR:HB	1:B:85:ILE:HG13	1.54	0.88
1:A:49:LEU:HD23	1:A:49:LEU:H	1.37	0.87
1:B:70:THR:HB	1:B:85:ILE:CG1	2.05	0.86
1:B:190:ARG:HH21	1:B:190:ARG:HG3	1.41	0.85
1:B:158:ASP:HB3	1:B:161:LYS:HG3	1.59	0.84
1:A:232:GLU:HG3	5:A:385:HOH:O	1.79	0.81
1:B:142:ILE:HD13	1:B:150:LYS:HG2	1.62	0.81
1:A:165:ASN:HD22	1:A:165:ASN:H	1.27	0.81
1:A:307:ALA:HB3	5:A:410:HOH:O	1.81	0.80
1:A:26:GLN:HE22	1:A:58:ASP:H	1.28	0.80

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:217:ILE:HA	5:A:391:HOH:O	1.81	0.79
1:B:50:ALA:O	1:B:54:HIS:HD2	1.66	0.78
1:A:81:LEU:HD13	1:A:82:GLU:N	1.99	0.76
1:B:44:ASN:HD22	1:B:58:ASP:HA	1.45	0.76
1:B:96:THR:O	1:B:98:PRO:HD3	1.86	0.75
1:A:18:ASP:HB3	5:A:367:HOH:O	1.87	0.75
1:A:240:ALA:HB1	1:A:248:TYR:HB3	1.69	0.75
1:A:136:PRO:HB2	5:A:419:HOH:O	1.88	0.74
1:B:142:ILE:HG21	1:B:150:LYS:HE2	1.68	0.74
1:B:224:ILE:HB	1:B:289:ILE:HD13	1.70	0.73
1:B:243:GLY:C	1:B:245:THR:H	1.91	0.73
1:B:136:PRO:HB2	1:B:139:TYR:HD1	1.53	0.72
1:B:103:ALA:HA	5:B:347:HOH:O	1.90	0.72
1:B:92:ILE:HG23	5:B:374:HOH:O	1.88	0.71
1:B:92:ILE:HG13	1:B:92:ILE:O	1.89	0.71
1:A:49:LEU:N	1:A:49:LEU:HD23	2.06	0.70
1:B:87:GLN:HE21	1:B:99:LYS:HA	1.56	0.70
1:A:115:PHE:HB3	5:A:425:HOH:O	1.92	0.69
1:B:43:SER:HB2	1:B:104:GLU:HB2	1.74	0.69
1:A:111:LEU:HD22	1:A:112:THR:HG23	1.74	0.69
1:B:164:GLU:O	1:B:165:ASN:HB3	1.93	0.69
1:A:165:ASN:N	1:A:165:ASN:HD22	1.84	0.68
1:B:70:THR:HA	5:B:400:HOH:O	1.94	0.67
1:A:160:SER:HB3	5:A:363:HOH:O	1.94	0.66
1:B:110:GLY:HA2	5:B:393:HOH:O	1.94	0.66
1:B:203:LEU:HD11	1:B:259:PRO:HG2	1.77	0.65
1:B:66:LYS:HB2	1:B:87:GLN:HB3	1.78	0.65
1:A:49:LEU:HG	5:A:416:HOH:O	1.95	0.65
1:B:81:LEU:HD13	1:B:81:LEU:C	2.18	0.65
1:B:133:LYS:HB3	3:B:330:NAG:H3	1.79	0.64
1:A:28:PHE:HE2	1:A:55:SER:O	1.80	0.64
1:A:50:ALA:O	1:A:54:HIS:HD2	1.80	0.64
1:A:43:SER:HB2	1:A:104:GLU:HB2	1.78	0.64
1:B:136:PRO:O	1:B:139:TYR:HB2	1.98	0.64
1:B:71:GLU:HG2	5:B:402:HOH:O	1.96	0.64
1:A:174:ILE:HB	5:A:409:HOH:O	1.97	0.63
1:B:164:GLU:O	1:B:165:ASN:CB	2.46	0.63
1:A:224:ILE:HG13	5:A:410:HOH:O	1.99	0.63
1:B:161:LYS:O	1:B:163:THR:N	2.32	0.62
1:B:87:GLN:HE21	1:B:99:LYS:CA	2.13	0.61
1:A:92:ILE:HD11	1:A:171:PHE:CE2	2.35	0.61
1:B:197:LYS:O	1:B:265:PHE:HA	2.01	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:90:LEU:HD11	1:B:92:ILE:HG22	1.83	0.60
1:A:255:ARG:NH2	5:A:370:HOH:O	2.34	0.60
1:A:40:TRP:CZ2	1:A:121:ILE:HG12	2.37	0.60
1:B:223:LEU:HB2	5:B:349:HOH:O	2.02	0.60
1:A:135:VAL:HG11	1:A:140:ASN:ND2	2.16	0.60
1:A:96:THR:O	1:A:98:PRO:HD3	2.02	0.59
1:A:163:THR:HA	5:A:414:HOH:O	2.03	0.59
1:B:137:PRO:HD3	5:B:417:HOH:O	2.00	0.59
1:B:65:TYR:CG	1:B:66:LYS:N	2.71	0.59
1:B:49:LEU:CD1	1:B:53:LEU:HD11	2.32	0.59
1:B:26:GLN:HE22	1:B:58:ASP:H	1.51	0.58
1:A:50:ALA:HA	1:A:53:LEU:HD12	1.83	0.58
1:A:19:ILE:HG22	1:A:92:ILE:HB	1.85	0.58
1:B:126:TYR:CD2	1:B:191:LYS:HB3	2.38	0.58
1:A:126:TYR:CD2	1:A:191:LYS:HB3	2.39	0.58
1:B:14:GLN:NE2	1:B:118:PHE:HD2	2.02	0.58
1:B:38:ASN:HD21	1:B:131:VAL:HB	1.69	0.58
1:B:88:ASP:OD1	1:B:89:THR:N	2.37	0.57
1:B:281:VAL:HG23	1:B:281:VAL:O	2.03	0.57
1:A:44:ASN:HA	1:A:56:LYS:HB3	1.87	0.57
1:A:26:GLN:NE2	1:A:58:ASP:H	2.00	0.57
1:A:314:SER:HB2	1:A:316:TYR:CE1	2.39	0.57
1:B:232:GLU:HG3	5:B:366:HOH:O	2.03	0.57
1:A:14:GLN:NE2	1:A:118:PHE:HD2	2.02	0.57
1:A:163:THR:O	1:A:164:GLU:C	2.42	0.57
1:A:182:ASP:N	1:A:182:ASP:OD1	2.38	0.57
1:B:203:LEU:HD12	1:B:260:ASP:O	2.05	0.56
1:B:9:ASN:ND2	1:B:159:THR:HG23	2.20	0.56
1:B:270:PHE:CZ	1:B:325:LEU:HD12	2.41	0.56
1:A:14:GLN:HE22	1:A:118:PHE:HD2	1.52	0.56
1:B:50:ALA:O	1:B:54:HIS:CD2	2.53	0.56
1:A:97:ILE:N	1:A:97:ILE:HD12	2.20	0.55
1:A:222:SER:O	1:A:223:LEU:HD23	2.06	0.55
1:A:240:ALA:HA	1:A:249:THR:O	2.06	0.55
1:A:175:ASP:OD1	1:A:177:SER:OG	2.24	0.55
1:B:70:THR:HB	1:B:85:ILE:HG12	1.83	0.55
1:B:70:THR:O	1:B:85:ILE:HG12	2.07	0.55
1:B:110:GLY:O	1:B:111:LEU:C	2.46	0.55
1:B:92:ILE:HD11	1:B:171:PHE:HE2	1.72	0.55
1:B:42:PRO:HB2	1:B:56:LYS:HG2	1.89	0.54
1:B:225:THR:HG23	5:B:411:HOH:O	2.08	0.54
1:A:198:PHE:HB2	1:A:215:ALA:HB2	1.88	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:73:ALA:HB2	1:A:82:GLU:HG3	1.90	0.54
1:A:65:TYR:HE1	1:A:86:SER:HB3	1.73	0.54
1:B:189:ARG:HB3	1:B:195:GLU:HG2	1.89	0.53
1:A:9:ASN:ND2	1:A:159:THR:HG23	2.23	0.53
1:B:142:ILE:HD13	1:B:150:LYS:CG	2.36	0.53
1:A:161:LYS:HB3	1:A:163:THR:HG23	1.91	0.53
1:B:11:LEU:C	1:B:13:ALA:H	2.12	0.53
1:A:3:HIS:HD2	1:A:171:PHE:O	1.90	0.53
1:B:174:ILE:HD12	1:B:179:PHE:CZ	2.44	0.53
1:B:100:GLN:OE1	1:B:140:ASN:ND2	2.42	0.53
1:A:306:ASP:HA	1:A:309:LEU:HB2	1.91	0.53
1:B:14:GLN:HE22	1:B:117:LYS:H	1.56	0.53
1:B:33:ASP:OD1	1:B:220:GLY:HA3	2.09	0.53
1:A:83:GLY:HA3	1:A:104:GLU:O	2.09	0.53
1:A:59:HIS:HD2	1:A:65:TYR:CZ	2.26	0.53
1:B:250:LEU:HD13	5:B:379:HOH:O	2.09	0.53
1:B:252:CYS:HA	5:B:379:HOH:O	2.08	0.52
1:A:307:ALA:HA	1:A:310:ARG:NH1	2.24	0.52
1:A:12:ASN:OD1	1:A:159:THR:HG23	2.10	0.52
1:B:155:TYR:HB2	1:B:313:TYR:HE1	1.74	0.52
1:B:3:HIS:CD2	1:B:95:LEU:HD13	2.45	0.52
1:B:33:ASP:HB3	5:B:382:HOH:O	2.09	0.52
1:A:85:ILE:HG22	1:A:86:SER:N	2.23	0.52
1:B:174:ILE:HG21	1:B:315:ILE:HD13	1.91	0.52
1:A:217:ILE:HG23	5:A:391:HOH:O	2.08	0.52
1:A:165:ASN:N	1:A:165:ASN:ND2	2.56	0.52
1:B:243:GLY:C	1:B:245:THR:N	2.60	0.52
1:A:203:LEU:HD12	1:A:260:ASP:O	2.09	0.52
1:A:187:PRO:HD2	1:A:266:ASN:ND2	2.25	0.52
1:B:301:LEU:CD1	5:B:375:HOH:O	2.58	0.51
1:B:285:CYS:C	1:B:286:ILE:HD12	2.31	0.51
1:B:24:PRO:HD2	1:B:61:ALA:O	2.11	0.51
1:A:59:HIS:HB2	1:A:65:TYR:CG	2.45	0.51
1:B:11:LEU:C	1:B:13:ALA:N	2.64	0.51
1:B:19:ILE:HG22	1:B:92:ILE:HB	1.93	0.51
1:B:46:CYS:HB2	1:B:105:ALA:O	2.11	0.51
1:A:175:ASP:C	1:A:175:ASP:OD1	2.48	0.51
1:A:311:LYS:O	1:A:311:LYS:CG	2.58	0.51
1:A:304:VAL:HB	5:A:391:HOH:O	2.10	0.51
1:B:54:HIS:HE1	1:B:116:GLY:O	1.93	0.51
1:A:75:GLN:CD	1:A:75:GLN:N	2.64	0.51
1:B:21:LEU:HD23	1:B:90:LEU:HA	1.93	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:161:LYS:HD3	5:B:392:HOH:O	2.10	0.50
1:B:123:GLY:HA2	5:B:382:HOH:O	2.10	0.50
1:B:160:SER:OG	1:B:160:SER:O	2.23	0.50
1:A:81:LEU:HD13	1:A:82:GLU:H	1.74	0.50
1:A:174:ILE:HG13	1:A:174:ILE:O	2.10	0.50
1:B:195:GLU:HG3	5:B:375:HOH:O	2.11	0.50
1:A:65:TYR:CG	1:A:66:LYS:N	2.80	0.50
1:A:49:LEU:CD2	1:A:49:LEU:H	2.03	0.50
1:A:151:ARG:NH2	5:A:345:HOH:O	2.42	0.50
1:B:20:THR:OG1	1:B:91:SER:HB2	2.11	0.50
1:B:190:ARG:HG3	1:B:190:ARG:NH2	2.19	0.50
1:B:282:SER:C	1:B:284:SER:H	2.15	0.50
1:B:195:GLU:CG	5:B:375:HOH:O	2.60	0.50
1:A:20:THR:OG1	1:A:91:SER:HB2	2.12	0.49
1:A:252:CYS:SG	1:A:280:GLU:HG3	2.51	0.49
1:B:85:ILE:HG22	1:B:86:SER:N	2.27	0.49
1:A:247:GLN:N	1:A:247:GLN:OE1	2.46	0.49
1:B:185:TRP:O	1:B:186:LEU:HD23	2.12	0.49
1:A:26:GLN:HE22	1:A:58:ASP:N	2.02	0.49
1:B:250:LEU:HD12	1:B:250:LEU:N	2.28	0.49
1:B:155:TYR:HB2	1:B:313:TYR:CE1	2.47	0.49
1:B:188:VAL:HG23	1:B:321:ASN:O	2.12	0.49
1:A:174:ILE:HD12	1:A:179:PHE:HE2	1.76	0.48
1:B:280:GLU:HG3	1:B:285:CYS:SG	2.53	0.48
1:A:198:PHE:O	1:A:213:HIS:HB2	2.14	0.48
1:B:222:SER:HA	1:B:306:ASP:HB2	1.95	0.48
1:A:81:LEU:CD1	1:A:82:GLU:N	2.76	0.48
1:B:87:GLN:NE2	1:B:99:LYS:CB	2.76	0.48
3:B:330:NAG:O7	3:B:331:NAG:O4	2.31	0.48
1:B:141:ALA:O	1:B:146:LEU:HB2	2.14	0.48
1:A:23:THR:O	1:A:62:SER:HA	2.14	0.48
1:B:131:VAL:O	1:B:134:VAL:HG23	2.13	0.47
1:A:46:CYS:HB2	1:A:105:ALA:O	2.14	0.47
1:A:299:GLY:HA2	1:A:300:PRO:C	2.34	0.47
1:A:200:GLY:N	1:A:264:ASN:HB3	2.29	0.47
1:B:316:TYR:CD1	1:B:316:TYR:N	2.81	0.47
1:A:261:LEU:HD23	1:A:263:PHE:CE1	2.49	0.47
1:B:48:SER:O	1:B:51:CYS:N	2.47	0.47
1:B:217:ILE:HG13	1:B:323:VAL:HG11	1.96	0.47
1:A:50:ALA:O	1:A:54:HIS:CD2	2.64	0.47
1:A:85:ILE:CG2	1:A:86:SER:N	2.78	0.47
1:A:87:GLN:NE2	1:A:99:LYS:HG2	2.29	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:175:ASP:OD2	1:B:178:LYS:HG2	2.14	0.46
1:A:143:GLN:HG3	1:A:144:GLN:HG3	1.97	0.46
1:B:59:HIS:HB2	1:B:65:TYR:CG	2.50	0.46
1:B:94:ASP:OD1	1:B:95:LEU:N	2.48	0.46
1:A:161:LYS:CB	1:A:163:THR:HG23	2.46	0.46
1:B:250:LEU:HD12	1:B:250:LEU:H	1.81	0.46
1:A:143:GLN:HG3	1:A:144:GLN:N	2.30	0.46
1:B:54:HIS:HB3	1:B:119:ASP:OD2	2.16	0.46
1:B:3:HIS:CD2	1:B:95:LEU:CD1	2.99	0.46
1:B:38:ASN:ND2	1:B:131:VAL:HB	2.32	0.45
1:B:294:PHE:HB3	1:B:298:VAL:CG1	2.45	0.45
1:A:65:TYR:OH	1:A:67:ALA:HA	2.16	0.45
1:B:94:ASP:OD1	1:B:94:ASP:C	2.53	0.45
1:B:225:THR:HA	1:B:290:THR:O	2.17	0.45
1:B:123:GLY:N	5:B:382:HOH:O	2.49	0.45
1:B:135:VAL:N	5:B:408:HOH:O	2.49	0.45
1:B:81:LEU:CD1	1:B:81:LEU:C	2.84	0.45
1:B:46:CYS:HB2	1:B:106:THR:HA	1.99	0.45
1:B:34:THR:OG1	1:B:218:ASP:HA	2.16	0.45
1:B:196:VAL:HB	1:B:266:ASN:HD21	1.82	0.45
1:B:153:ALA:O	1:B:170:THR:N	2.50	0.45
1:A:65:TYR:CZ	1:A:67:ALA:HA	2.52	0.45
1:A:66:LYS:HB2	1:A:87:GLN:HB3	1.98	0.45
1:B:51:CYS:O	1:B:52:PHE:C	2.55	0.45
1:B:190:ARG:NH1	1:B:298:VAL:O	2.51	0.45
1:B:168:GLU:HG2	1:B:169:ALA:O	2.17	0.45
1:A:198:PHE:HB3	1:A:213:HIS:HB3	1.98	0.44
1:A:87:GLN:HE22	1:A:99:LYS:HG2	1.82	0.44
1:A:256:ASP:C	1:A:258:LEU:H	2.20	0.44
1:A:311:LYS:O	1:A:311:LYS:HG2	2.16	0.44
1:B:65:TYR:OH	1:B:67:ALA:HA	2.16	0.44
1:B:85:ILE:HA	5:B:347:HOH:O	2.17	0.44
1:A:72:PHE:CD2	1:A:72:PHE:C	2.91	0.44
1:B:23:THR:O	1:B:62:SER:HA	2.17	0.44
1:B:40:TRP:NE1	1:B:121:ILE:HB	2.33	0.44
1:B:92:ILE:HD11	1:B:171:PHE:CE2	2.52	0.44
1:B:26:GLN:NE2	1:B:58:ASP:H	2.14	0.44
1:A:44:ASN:CB	1:A:57:TYR:O	2.66	0.44
1:A:87:GLN:HG2	1:A:88:ASP:N	2.32	0.44
1:B:150:LYS:O	1:B:151:ARG:HB3	2.18	0.43
1:B:11:LEU:O	1:B:13:ALA:N	2.51	0.43
1:A:59:HIS:N	1:A:59:HIS:ND1	2.64	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:151:ARG:HB2	1:B:317:ASP:HA	2.00	0.43
1:A:44:ASN:HB3	1:A:57:TYR:O	2.19	0.43
1:A:39:LEU:HD13	1:A:122:LEU:HG	2.00	0.43
1:A:138:PHE:CE2	1:A:142:ILE:HD11	2.52	0.43
1:B:134:VAL:HG13	3:B:331:NAG:O6	2.19	0.43
1:B:29:LYS:O	1:B:119:ASP:N	2.44	0.43
1:B:87:GLN:NE2	1:B:99:LYS:HG2	2.33	0.43
1:B:282:SER:O	1:B:284:SER:N	2.47	0.43
1:B:2:GLY:HA2	5:B:362:HOH:O	2.17	0.43
1:A:14:GLN:NE2	1:A:118:PHE:CD2	2.85	0.43
1:A:9:ASN:HD21	1:A:159:THR:HG23	1.83	0.43
1:A:151:ARG:CB	1:A:317:ASP:HA	2.48	0.43
1:A:315:ILE:O	1:A:323:VAL:HA	2.18	0.43
1:B:31:ILE:HB	1:B:121:ILE:HD13	2.00	0.43
1:A:85:ILE:HD13	1:A:103:ALA:HB2	2.01	0.43
1:B:299:GLY:HA2	1:B:300:PRO:C	2.39	0.43
1:B:65:TYR:HE1	1:B:86:SER:HB3	1.84	0.43
1:B:155:TYR:CE1	1:B:329:ILE:HG13	2.54	0.43
1:B:134:VAL:HG22	3:B:331:NAG:C6	2.49	0.42
1:A:40:TRP:NE1	1:A:121:ILE:HB	2.33	0.42
1:A:163:THR:HG21	5:A:349:HOH:O	2.17	0.42
1:A:85:ILE:CD1	1:A:103:ALA:HB2	2.49	0.42
1:A:100:GLN:OE1	1:A:137:PRO:HA	2.19	0.42
1:B:178:LYS:O	1:B:179:PHE:HB3	2.18	0.42
1:B:157:GLY:N	1:B:166:GLY:O	2.42	0.42
1:A:87:GLN:O	1:A:88:ASP:HB2	2.19	0.42
1:B:286:ILE:HD12	1:B:286:ILE:N	2.34	0.42
1:A:311:LYS:HG3	1:A:327:LYS:HE2	2.00	0.42
1:B:8:THR:HB	1:B:16:TYR:CE2	2.54	0.42
1:B:272:ILE:HB	1:B:276:ASP:HB2	2.02	0.42
1:A:3:HIS:CD2	1:A:171:PHE:O	2.70	0.42
1:B:195:GLU:HA	1:B:215:ALA:O	2.19	0.42
1:A:315:ILE:O	1:A:324:GLY:N	2.38	0.42
1:A:207:TYR:CD1	1:A:208:ALA:N	2.87	0.42
1:A:11:LEU:C	1:A:13:ALA:N	2.72	0.42
1:B:135:VAL:HA	1:B:136:PRO:HD3	1.85	0.42
1:A:316:TYR:N	1:A:316:TYR:CD1	2.87	0.42
1:A:28:PHE:CE2	1:A:55:SER:O	2.68	0.42
1:B:58:ASP:OD1	1:B:60:GLU:OE2	2.38	0.42
1:A:251:ASP:OD2	1:A:251:ASP:C	2.58	0.42
1:A:49:LEU:N	1:A:49:LEU:CD2	2.73	0.41
1:A:309:LEU:HD23	1:A:309:LEU:HA	1.79	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:238:ILE:HB	1:A:258:LEU:HD13	2.02	0.41
1:A:188:VAL:HG23	1:A:321:ASN:O	2.19	0.41
1:B:68:ASN:OD1	1:B:70:THR:OG1	2.20	0.41
1:B:174:ILE:HD12	1:B:179:PHE:CE2	2.56	0.41
1:B:136:PRO:HD2	1:B:139:TYR:CD1	2.55	0.41
1:A:43:SER:C	1:A:45:GLU:H	2.23	0.41
1:A:161:LYS:O	1:A:162:ASP:HB2	2.20	0.41
1:A:21:LEU:HD23	1:A:90:LEU:HA	2.03	0.41
1:B:178:LYS:HD2	1:B:329:ILE:HD11	2.02	0.41
1:B:179:PHE:CD2	1:B:313:TYR:CD2	3.09	0.41
1:B:320:ASN:O	1:B:321:ASN:C	2.59	0.41
1:B:248:TYR:CD1	1:B:248:TYR:N	2.88	0.41
1:A:276:ASP:O	1:A:307:ALA:HB1	2.21	0.41
1:A:59:HIS:CD2	1:A:65:TYR:CZ	3.08	0.41
1:B:9:ASN:HD21	1:B:159:THR:HG23	1.85	0.41
1:B:34:THR:HG1	1:B:218:ASP:HA	1.85	0.41
1:B:318:LEU:HD12	1:B:318:LEU:O	2.20	0.41
1:B:87:GLN:NE2	1:B:99:LYS:CA	2.82	0.41
1:B:8:THR:HB	1:B:16:TYR:CD2	2.55	0.41
1:A:329:ILE:OXT	1:A:329:ILE:HG13	2.21	0.41
1:A:189:ARG:HH21	1:A:214:GLY:CA	2.34	0.41
1:B:5:VAL:HG21	1:B:92:ILE:CD1	2.34	0.40
1:B:59:HIS:CB	1:B:65:TYR:CD1	3.04	0.40
1:A:307:ALA:CB	5:A:410:HOH:O	2.55	0.40
1:A:174:ILE:HD12	1:A:179:PHE:CE2	2.56	0.40
1:B:198:PHE:HB3	1:B:213:HIS:HB3	2.04	0.40
1:B:86:SER:N	5:B:347:HOH:O	2.55	0.40
1:A:250:LEU:N	1:A:250:LEU:CD1	2.83	0.40
1:A:27:ASN:HB3	5:A:367:HOH:O	2.21	0.40
1:B:144:GLN:O	1:B:145:ASP:C	2.60	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	321/329 (98%)	297 (92%)	20 (6%)	4 (1%)	19	38
1	B	321/329 (98%)	282 (88%)	24 (8%)	15 (5%)	4	4
All	All	642/658 (98%)	579 (90%)	44 (7%)	19 (3%)	7	10

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	109	PRO
1	A	111	LEU
1	B	111	LEU
1	B	162	ASP
1	B	163	THR
1	B	165	ASN
1	B	176	GLU
1	B	145	ASP
1	A	164	GLU
1	A	192	ALA
1	B	112	THR
1	B	160	SER
1	B	321	ASN
1	B	12	ASN
1	B	133	LYS
1	B	144	GLN
1	B	192	ALA
1	B	283	GLY
1	B	300	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	265/268 (99%)	241 (91%)	24 (9%)	14	25
1	B	265/268 (99%)	240 (91%)	25 (9%)	13	23
All	All	530/536 (99%)	481 (91%)	49 (9%)	13	24

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	VAL
1	A	6	PRO
1	A	49	LEU
1	A	59	HIS
1	A	81	LEU
1	A	82	GLU
1	A	92	ILE
1	A	101	ASP
1	A	111	LEU
1	A	143	GLN
1	A	162	ASP
1	A	165	ASN
1	A	180	LYS
1	A	201	ILE
1	A	205	ASP
1	A	230	LEU
1	A	245	THR
1	A	250	LEU
1	A	256	ASP
1	A	279	LEU
1	A	281	VAL
1	A	298	VAL
1	A	308	PHE
1	A	314	SER
1	B	5	VAL
1	B	18	ASP
1	B	49	LEU
1	B	59	HIS
1	B	90	LEU
1	B	92	ILE
1	B	104	GLU
1	B	117	LYS
1	B	127	ASP
1	B	137	PRO
1	B	143	GLN
1	B	145	ASP
1	B	160	SER
1	B	163	THR
1	B	180	LYS
1	B	190	ARG
1	B	205	ASP
1	B	227	PRO
1	B	228	SER

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Mol	Chain	Res	Type
1	B	230	LEU
1	B	245	THR
1	B	250	LEU
1	B	271	THR
1	B	279	LEU
1	B	327	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	HIS
1	A	14	GLN
1	A	26	GLN
1	A	54	HIS
1	A	87	GLN
1	A	140	ASN
1	A	143	GLN
1	A	165	ASN
1	A	235	ASN
1	A	266	ASN
1	B	3	HIS
1	B	14	GLN
1	B	54	HIS
1	B	87	GLN
1	B	235	ASN
1	B	266	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

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	$\# Z > 2$
2	BMA	A	330	2	10,11,12	0.45	0	11,15,17	0.41	0
2	NAG	A	331	2	12,14,15	0.68	0	15,19,21	0.85	1 (6%)
2	NAG	A	332	1,2	12,14,15	0.51	0	15,19,21	0.97	1 (6%)
3	NAG	B	330	3	12,14,15	0.52	0	15,19,21	0.75	0
3	NAG	B	331	1,3	12,14,15	0.67	0	15,19,21	1.18	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
2	BMA	A	330	2	-	0/2/19/22	0/1/1/1
2	NAG	A	331	2	-	0/6/23/26	0/1/1/1
2	NAG	A	332	1,2	-	0/6/23/26	0/1/1/1
3	NAG	B	330	3	-	0/6/23/26	0/1/1/1
3	NAG	B	331	1,3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	331	NAG	C3-C4-C5	2.56	114.78	110.20
2	A	331	NAG	C3-C2-N2	-2.24	108.35	111.76
3	B	331	NAG	C3-C2-N2	-2.12	108.54	111.76
2	A	332	NAG	C6-C5-C4	-2.08	107.96	113.00

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry

2 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	$\# Z > 2$
4	NAG	A	333	1	12,14,15	0.78	0	15,19,21	0.93	1 (6%)
4	NAG	B	332	1	12,14,15	0.57	0	15,19,21	0.95	1 (6%)

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	A	333	1	-	0/6/23/26	0/1/1/1
4	NAG	B	332	1	-	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	A	333	NAG	C2-N2-C7	-2.55	118.81	123.09
4	B	332	NAG	C2-N2-C7	-2.33	119.17	123.09

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	325/329 (98%)	-0.28	7 (2%) 59 56	26, 51, 82, 114	31 (9%)
1	B	324/329 (98%)	-0.29	8 (2%) 54 52	25, 54, 82, 103	37 (11%)
All	All	649/658 (98%)	-0.28	15 (2%) 57 55	25, 52, 82, 114	68 (10%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	245	THR	9.8
1	B	11	LEU	3.5
1	B	244	TRP	3.4
1	B	164	GLU	3.1
1	A	246	GLY	2.9
1	B	39	LEU	2.8
1	A	76	TYR	2.8
1	B	246	GLY	2.7
1	A	164	GLU	2.7
1	A	47	GLY	2.6
1	A	1	GLY	2.4
1	B	22	GLY	2.3
1	B	127	ASP	2.1
1	A	162	ASP	2.1
1	B	63	SER	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 ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	NAG	A	332	14/15	0.29	2.63	42,62,76,77	0
3	NAG	B	330	14/15	0.28	1.94	46,58,76,91	0
3	NAG	B	331	14/15	0.21	1.47	58,82,96,99	0
2	NAG	A	331	14/15	0.12	-0.88	48,58,61,65	2
2	BMA	A	330	11/12	0.14	-7.67	58,65,68,71	2

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	NAG	B	332	14/15	0.17	1.30	48,60,66,70	0
4	NAG	A	333	14/15	0.17	-0.08	32,52,64,66	2

6.5 Other polymers ⓘ

There are no such residues in this entry.